The listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims

Claim 1 (canceled).

Claim 2 (currently amended): The method of Claim 14, wherein R¹⁵ is selected from H, optionally substituted pyrrolidinyl, optionally substituted piperazinyl, optionally substituted piperidinyl, morpholinyl, 1,2,3,6-tetrahydro-pyridinyl, (optionally substituted pyrrolidinyl)-C1-C2-alkyl, (optionally substituted piperidinyl)-C1-C2-alkyl, (optionally substituted piperazinyl)-C1-G2-alkyl, morpholinyl-G1-G2-alkyl, G1-G4-alkylamino-G1-G4-alkyl, G1-G4-hydroxyalkylamino, (eptionally substituted pyrrelidinyl)-C1-C2-alkylamine, (eptionally substituted piperidinyl)-C1-C2-alkylamino, (optionally substituted piperazinyl)-C1-C2-alkylamino, morpholinyl-C1-C2alkylamino, optionally substituted pyrrolidinyl-C₁-C₄-alkoxy, optionally substituted azetidinyl-C₁-C₄-alkoxy, tetrahydrofuryl-C₁-C₄-alkoxy, optionally substituted piperidinyl-C₁-C₄-alkoxy, C₁-C₄-alkylamino-C₁-C₄-alkoxy, tetrahydrofuryloxy, optionally substituted piperidinyloxy, and optionally substituted phenoxy, G₄-G₄-alkylaminocarbonyl and G₄-G₄-alkylaminothiocarbonyl; wherein R¹⁶ is selected from H. 5-6-membered nitrogen-containing heterocyclylearbonyl, C₁-C₄alkylaminocarbonyl, and C₁-C₄-alkylaminomethyl, and 5-6-membered nitrogen containing heteroeyelylmethyl; and wherein R¹⁷ is selected from halo, C₁-C₂-alkyl, thienylsulfonyl- C₁-C₂alkyl, optionally substituted 5-6-membered heteroarylsulfonyl-C4-C2-alkyl, optionally substituted phenoxy: and C₃-C₆-cycloalkyl-C₂-C₄-alkynyl: and pharmaceutically acceptable derivatives thereof.

Claim 3 (currently amended): The method of Claim 2, wherein R¹⁵ is selected from H, tetrahydro-furanyloxy, 1-methylpyrrolidin-2-ylmethoxy, 2-pyrrolidinylmethoxy, 3-pyrrolidinylmethoxy, 1-Boc-pyrrolidin-2-ylmethoxy, 4-piperidinylmethoxy, 1-Boc-piperidin-4-ylethoxy, piperidin-4-ylethoxy, 1-methyl-piperidin-4-ylmethoxy, 1-Boc-azetidin-3-ylmethoxy, 1-methyl-azetidin-3-ylmethoxy, 3-azetidinylmethoxy, 1-methyl-piperidin-4-yloxy, phenoxy, 4-(pyrrolidin-1-ylmethyl)phenoxy, and dimethylaminoethoxy, 1-piperidinylmethyl, 1-(piperidin-1-yl)ethyl, 3-methylpiperidin-1-yl)ethyl, 3-methylpiperidin-1-yl

ylmethyl, 1-pyrrolidinylmethyl, 2,2,6,6-tetramethylpiperidin-1-ylmethyl, 2,6-dimethylpiperidin-1-ylmethyl, diethylaminomethyl, diethylaminomethyl, diethylaminomethyl, diethylaminomethyl, diethylaminomethyl, 2-thienylsulfonylmethyl, hydroxypropylamino, 4-ethyl piperidin-1-yl, 4-(2-pyridyl)piperidin-1-yl, 1-methyl-1,2,3,6-tetrahydro-pyridin-4-yl, 1-methyl-1,2,3,6-tetrahydro-pyridin-4-yl, 1-2,3,6-tetrahydro-pyridin-4-yl, wherein R¹⁶ is selected from H, 1-piperidinylearbonyl, diethylaminocarbonyl, and diethylaminomethyl, 1-piperidinylmethyl; and wherein R¹⁷ is selected from chloro, bromo, methyl and cyclopropylethynyl, and pharmaccutically acceptable derivatives thereof.

Claim 4 (currently amended): The method of Claim 3, wherein R¹⁷ is chloro or bromo; and pharmaceutically acceptable derivatives thereof.

Claim 5 (currently amended): The method of Claim 14, wherein R 15 is selected from H, optionally substituted pyrrolidinyl, optionally substituted piperazinyl, optionally substituted piperazinyl, optionally substituted piperazinyl, (optionally substituted pyrrolidinyl) C₁-C₂-alkyl, (optionally substituted piperazinyl) C₄-C₅-alkyl, (optionally substituted piperazinyl) C₄-C₅-alkyl, morpholinyl-C₄-C₅-alkyl, (optionally substituted pyrrolidinyl) C₄-C₅-alkylamino, (optionally substituted piperazinyl) C₄-C₅-alkylamino, morpholinyl-C₄-C₅-alkylamino, C₄-C₄-alkylamino-C₄-C₄-alkyl, C₄-C₄-alkylamino, optionally substituted pyrrolidinyl-C₁-C₄-alkoxy, optionally substituted azetidinyl-C₁-C₄-alkoxy, tetrahydrofuryl-C₁-C₄-alkoxy, optionally substituted piperidinyl-C₁-C₄-alkoxy, ctrahydrofuryloxy, optionally substituted piperidinyloxy, and optionally substituted phenoxy, C₄-C₄-alkylaminocarbonyl and C₄-C₄-alkylaminothiocarbonyl; wherein R 16 is selected from H, 5-6-membered nitrogen-containing heterocyclylmethyl; and wherein R 17 is selected from C₃-C₆-cycloalkyl and phenyl-optionally substituted with one or two substituents selected from halo,

C₁-C₄-alkylamino, amino, nitro, C₁-C₄-alkoxy, C₁-C₂-haloalkyl, hydroxy, C₁-C₄-alkylthio, C₄-C₄-alkylcarbonylamino, (optionally substituted phenyl)sulfonylamino, eyano, C₄-C₂-haloalkoxy, 5-or 6-membered N-containing heterocyclyl, aminosulfonyl,

(6-membered N-containing heterocyclyl)sulfonyl, G₄- G₂-haloalkylearbonylaminosulfonyl and (optionally substituted-phenyl)aminosulfonyl; and phermacoutically acceptable derivatives thereof.

Claim 6 (currently amended): The method of Claim 5, wherein R¹⁵ is selected from H, tetrahydro-furanyloxy, 1-methylpyrrolidin-2-vlmethoxy, 2-pyrrolidinylmethoxy, 3pyrrolidinylmethoxy, 1-Boc-pyrrolidin-2-ylmethoxy, 4-piperidinylmethoxy, 1-Boc-piperidin-4vlmethoxy, 1-Boc-piperidin-4-ylethoxy, piperidin-4-ylethoxy, 1-methyl-piperidin-4-ylmethoxy, 1-Boc-azetidin-3-ylmethoxy, 1-methyl-azetidin-3-ylmethoxy, 3-azetidinylmethoxy, 1-methylpiperidin-4-yloxy, phenoxy 4-(pyrrolidin-1-ylmethyl)phenoxy, and dimethylaminoethoxy, 1-piperidinylmethyl, 1-(piperidin-1-yl)ethyl, 3-methylpiperidin-1vimethyl, 1-pyrrolidinylmethyl, 2,2,6,6-tetramethylpiporidin-1-ylmethyl, 2,6-dimethylpiporidin-1-vlmethyl.-dimethylaminomethyl. diethylaminomethyl, diethylaminothiocarbonyl; diethylaminocarbonyl, N-Boe-N-isopropylaminomethyl, isopropylaminomethyl, 2thienvisulfonvimethyl, hydroxypropylamino, 4 ethyl-piperidin 1-yl, 4 (2-pyridyl)piperidin 1-yl, 1-mothylpiperidin 4 vl. 4 (2-pyrazinyl)piperidin 1-yl. 1-methyl-1,2,3,6-tetrahydro-pyridin 4-yl, 1,2,3,6 tetrahydro-pyridin 4 yl, and 1-Boe-1,2,3,6 tetrahydro-pyridin 4 yl; wherein R¹⁶ is selected from H. 1-piperidinylearbonyl, diethylaminocarbonyl, and diethylaminomethyl, 1piperidinylmethyl: and wherein R¹⁷ is selected from cyclopropyl and phenyl-optionally substituted with aminosulfonyl; and pharmaceutically acceptable derivatives thereof.

Claim 7 (currently amended): The method of Claim 6, wherein R¹⁷ is unsubstituted phenyl; and pharmaceutically acceptable derivatives thereof.

Claim 8 (currently amended): The method of Claim 14, wherein R¹⁵ is selected from H₃ optionally substituted pyrrolidinyl, optionally substituted piperazinyl, optionally substituted piperazinyl, optionally substituted pyrrolidinyl) C₄-C₂-alkyl, (optionally substituted piperazinyl) C₄-C₂-alkyl, morpholinyl C₄-C₂-alkyl, (optionally substituted piperazinyl) C₄-C₂-alkyl, morpholinyl C₄-C₂-alkyl, (optionally substituted pyrrolidinyl) C₄-C₂-alkylamine, (optionally substituted piperazinyl) C₄-C₂-alkylamine, morpholinyl C₄-C₂-alkylamine, C₄-alkylamine C₄-C₄-alkyl₄-C₄-alkyl₅-C₄-alkyl₆

azetidinyl-C₁-C₄-alkoxy, tetrahydrofuryl-C₁-C₄-alkoxy, optionally substituted piperidinyl-C₁-C₄-alkoxy, C₁-C₄-alkylamino-C₁-C₄-alkoxy, tetrahydrofuryloxy, optionally substituted piperidinyloxy, and optionally substituted phenoxy, C₄-C₄-alkylaminocarbonyl and C₁-C₄-alkylaminothiocarbonyl; wherein R¹⁶ is selected from H, 5-6-membered nitrogen-containing heterocyclylearbonyl, C₁-C₄-alkylaminocarbonyl, and C₁-C₄-alkylaminomethyl, and 5-6-membered nitrogen-containing heterocyclylmethyl; and wherein R¹⁷ is selected from optionally substituted indazolyl, optionally substituted indolyl, unsubstituted 5-membered oxygen or sulfur containing heterocyclyl, unsubstituted thienyl, unsubstituted 6-membered nitrogen-containing heterocyclyl, and 6-membered nitrogen-containing heterocyclyl substituted with one-or-more substituents independently selected from pyridyl, phenyl,

G₄-G₄-alkyl, G₄-G₂-haloalkyl, C₁-C₂ alkoxy, amino, halo, piperidinyl, morpholinyl, G₄-G₂ alkylaminothiocarbonyl, N,N-di-C₄-G₂ alkylamino-G₄-G₄-alkylenyl, N-G₄-G₂-alkylamino-G₄-G₄-alkylenyl, morpholinyl-G₄-G₄-alkylenylaminocarbonyl, aminocarbonyl, G₄-G₂-haloalkylearbonylamino, morpholinyl-G₄-G₄-alkylonylamino, N,N-di-G₄-G₂-alkylamino and N,N-di-G₄-G₂-alkylamino-G₄-G₄-alkylenylamino,

and pharmacoutically acceptable derivatives thereof.

Claim 9 (currently amended): The method of Claim 8, wherein R¹⁵ is selected from H, tetrahydro-furanyloxy, 1-methylpyrrolidin-2-ylmethoxy, 2-pyrrolidinylmethoxy, 3-pyrrolidinylmethoxy, 1-Boc-piperidin-4-ylmethoxy, 1-Boc-piperidin-4-ylmethoxy, 1-Boc-piperidin-4-ylmethoxy, 1-methyl-piperidin-4-ylmethoxy, 1-methyl-piperidin-4-ylmethoxy, 1-methyl-piperidin-4-ylmethoxy, 1-methyl-piperidin-4-yloxy, phenyloxy, 4-(pyrrolidin-1-ylmethyl)phenoxy, and dimethylaminoethoxy, 1-piperidinylmethyl, 1-(piperidin-1-yl)ethyl, 3-methylpiperidin-1-ylmethyl, 1-pyrrolidinylmethyl, 2,2,6,6 tetramethylpiperidin-1-ylmethyl, 2,6 dimethylpiperidin-1-ylmethyl, diethylaminomethyl, diethylaminomethyl, diethylaminomethyl, diethylaminomethyl, diethylaminomethyl, byl, 1-pyrrolidin-1-yl, 1-methyl-1,2,3,6 tetrahydro-pyridin-1-yl, 1-methyl-1,2,3,6 tetrahydro-pyridin-1-yl, 1-piperidin-1-yl, 1-piperid

wherein R¹⁷ is selected from 5 indazelyl; 1 Boe indel-5 yl; unsubstituted thienyl; 5 tert-butyloxazel 2 yl and 4-pyridyl substituted with-one or more substituents independently selected from methoxy and chloro; and pharmaceutically acceptable derivatives thereof.

Claim 10 (currently amended): The method of Claim 9 8, wherein R¹⁷ is 4-pyridyl; and pharmaceutically acceptable derivatives thereof.

Claim 11 (currently amended): The method of Claim 14 and pharmaceutically acceptable derivatives thereof, wherein the compound is selected from:

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1-[6-(3 Methyl piperidin 1 ylmethyl) pyridin 2 yl] 3 (2 phenyl thiazel 4 yl) urea;
1-[4 (Piperidine-1-earbonyl) pyridin-2-yl]-3-(2-pyridin-4-yl thiazol-4-yl) urea;
1-(2 Chloro-thiazol 4 yl) 3-[4-(piperidine-1-earbonyl) pyridin 2-yl] urea;
N.N. Diethyl 2-[3-(2 pyridin 4 yl thiazol 4 yl) ureido] isonicotinamide;
N.N. Diethyl 2-13-(2-phonyl thiazol-4-yl) urcidol-isonicotinamide;
2-13-(2-Bromo-thiazol-4-vl) urcidol-N.N-diethyl-isonicotinamide;
1 (4-Diethylaminomethyl-pyridin-2-yl) 3 (2-pyridin-4-yl-thiazol-4-yl)-urea;
1-[6 (2,6-Dimethyl-piperidin-1-ylmethyl) pyridin-2-yl] 3 (2-pyridin-4-yl-thiazol-4-yl) urea;
1-[6-(1-Piperidin-1-yl-ethyl)-pyridin-2-yl] 3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
2-(16-[3-(2-Pyridin-4-yl-thiazol-4-yl)-urcido]-pyridin-2-ylamino}-methyl)-piperidine-1-
     carboxylic acid tert-butyl ester;
1-{6-[(Piperidin-2-ylmethyl)-amino]-pyridin-2-yl}-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
(S) 1-[6-(3-Methyl-piperidin-1-ylmethyl) pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
(R) 1-[6-(3-Methyl-piperidin-1-ylmethyl)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
1-(2 Chloro-thiazol 4 yl) 3 (6 piperidin-1-ylmethyl-pyridin-2 yl) urea;
1-(2-Bromo-thiazol-4-yl)-3-[6-(2-piperidin-4-yl-ethoxy)-pyridin-2-yl]-urea;
1-(2-Chloro-thiazol-4-yl)-3-[6-(2-piperidin-4-yl-ethoxy)-pyridin-2-yl]-urea;
1-[6-(Azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-bromo-thiazol-4-yl)-urea;
1-[6-(Azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-chloro-thiazol-4-yl)-urea;
1-(2-Bromo-thiazol-4-yl)-3-[6-(piperidin-4-ylmethoxy)-pyridin-2-yl]-urea;
1-(2-Chloro-thiazol-4-yl)-3-[6-(piperidin-4-ylmethoxy)-pyridin-2-yl]-urea;
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tert-Butyl 3-{6-[3-(2-pyridin-4-yl-thiazol-4-yl)-ureido]-pyridin-2-yloxymethyl}-pyrrolidine-1-
     carboxylate;
1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(pyrrolidin-3-ylmethoxy)-pyridin-2-yl]-urea;
1-(2-Cyclopropyl-thiazol-4-yl)-3-[6-(2-piperidin-4-yl-ethoxy)-pyridin-2-yl]-urea;
1-[6 (Isopropylamine methyl) pyridin 2 yl] 3 (2 pyridin 4 yl thiazol 4 yl) urea;
1-[6 (Isopropylamino methyl) pyridin 2-yl] 3-(2-phenyl-thiazol-4-yl) urea;
1-(2-Bromo thiazol 4-yl)-3-[6-(isopropylamino-methyl)-pyridin-2-yl]-urea;
1-(2-Bromo-thiazol-4-yl)-3-[6-(1-methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-urea;
1-(2-Chloro-thiazol-4-yl)-3-[6-(1-methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-urea;
1-(2-phenylthiazol-4-yl)-3-(6-p-pyrrolidin-1-ylmethylphenoxypyridin-2-yl)urea;
1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(tetrahydro-furan-3-yloxy)-pyridin-2-yl]-urea;
1-[2-(111-Indazol 5-yl) thiazol 1-yl] 3 (6-piperidin-1-ylmethyl-pyridin 2-yl) urea;
1-(1' Methyl-1',2',3',6' tetrahydro [2,4']bipyridinyl-6-yl)-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
1-(2-Bromo thizol 4-yl) 3-(1'-methyl-1',2',3',6'-tetrahydro-[2,4']bipyridinyl-6-yl)-urca;
1-(1' Methyl 1',2',3',6'-tetrahydro-2[2,4]bipyridinyl-6-yl) 3-(2 phenyl-thiazol 4 yl) urea;
1-[6-(3 Hydroxy propylamino) pyridin 2-yl-3-(2-pyridin 4-yl-thizol 4-yl)-urca;
1-(2 Bromo thiazol 4 yl) 3 [6(3-hydroxy propylamine) pyridin 2 yl] urea;
1 (1' Methyl 1',2',3',4',5',6' hexahydro [2,4']bipydrinyl 6 yl) 3 (2 pyridin 4 yl thiazel 4
     uroa:
1-(1' Methyl-1',2',3',4',5',6'-hexahydro-[2,4']bipyridinyl-6-yl)-3-(2-phonyl-thiazol-4-yl)-urea;
6-[3-(2-Pyridin 4-yl-thizol-4-yl)-urcido]-3',6'-dihydro-2'H-[2,4]bipyridinyl-1'-carboxylic acid
     tert-butylester;
1-(2-Pyridin 4 vl thiazol 4 vl) 3 (1'2'3'.6' tetrahydro [2,4']bipyridinyl 6 vl) urca;
1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(tetrahydro-furan-3-ylmethoxy)-pyridin-2-yl]-urea;
2-[6-[3-(2-Pyridin-4-yl-thiazol-4-yl)-ureido]-pyridin-2-yloxymethyl]-pyrrolidine-1-carboxylic
     acid tert-butyl ester;
1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-urea;
6-[3-(2-Pyridin-4-yl-thiazol-4-yl)-ureido]-pyridino-2-carbothioic acid-dicthylamide;
1-(2-Bromo thiazol 4 yl) 3-[6 (3-methyl-piperidin-1 ylmethyl)-pyridin-2-yl]-urea;
1-(2 Chloro-thiazel 4 yl) 3-[6 (3 methyl-piperidin-1 ylmethyl) pyridin-2-yl] ureat
1-(2-Phonyl thiazol-4-yl)-3-[4 (piperidine-1-earbonyl) pyridin-2-yl]-urea;
1-(2-Bromo-thiazol 4 yl) 3-[4 (piperidine 1 carbonyl) pyridin-2-yl]-urea;
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1-[2-(2-Methoxy-pyridin-4-yl)-thiazol-4-yl]-3-(6-phenoxy-pyridin-2-yl)-urea;
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- 1-[2-(2-Methoxy-pyridin-4-yl)-thiazol-4-yl]-3-[6-(1-methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-urea;
- 1-[6-(2-Dimethylamino-ethoxy)-pyridin-2-yl]-3-[2-(2-methoxy-pyridin-4-yl)-thiazol-4-yl]-urea;
- 1-[6-(1-Methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-(2-phenylthiazol 4-yl) 3-(6-pyrrolidin-1-ylmethyl-pyridin-2-yl)urea;
- 1-(6-Diothylaminomethylpyridin-2-yl)-3-(2-phenylthiazol-4-yl)urca;
- (S)-1-[6-(1-Methylpyrrolidin-2-ylmethoxy)pyridin-2-yl]-3-(2-phenylthiazol-4-yl)urea;
- 1-[6-(2-Piperidin-4-yl-ethoxy)pyridin-2-yl]-3-[2-phenylthiazol-4-yl]urea;
- 1-[6-(4-Ethylpiperazin-1-yl)-pyridin-2-yl]-3-(2-phenylthiazol-4-yl)urea;

Diethyl 6-[3 (2-phonylthiazol 4-yl)uroido]-pyridine-2-carboxamide;

- 1-(2-Pyridin-4-yl-thiazol-4-yl)-3-(6-p-pyrrolidin-1-ylmethylphenoxypyridin-2-yl)urea;
- 1-(2-Bromothiazol-4-yl)-3-(6-p-pyrrolidin-1-ylmethylphenoxypyridin-2-yl)urea;
- 1-[6-(Piperidin-4-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-piperidin-4-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-piperidin-4-yloxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-[6-(Azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;
- 1-(2-Phenyl-thiazol-4-yl)-3-[6-(piperidin-4-ylmethoxy)-pyridin-2-yl]-urea;
- 1-[6-(1-Methyl-piperidin-4-ylmethoxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-piperidin-4-yloxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;
- 1-[6-(2-Piperidin-4-yl-ethoxy)-pyridin-2-yl]-3-(2-thiophen-2-yl-thiazol-4-yl)-urea; and
- 1-[6-(1-Methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-3-[2-(thiophene-2-sulfonylmethyl)-thiazol-4-yl]-urea;
- 1-[2 (2-Methoxy-pyridin 4-yl) thiazol 4-yl] 3 (6-piperdin 1-ylmethyl-pyridin 2-yl) urea; and [2 (2-Chloro-pyridin 4-yl) thiazol 4-yl] 3 (6-piperdin 1-ylmethyl-pyridin 2-yl) urea; and pharmaceutically acceptable salts thereof.
- Claim 12 (currently amended): The method of Claim 14 and pharmaccutically acceptable derivatives thereof, wherein the compound is selected from:

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1-[6-(3-Methyl-piperidin-1-ylmethyl) pyridin-2-yl]-3-(2-phenyl-thiazel-4-yl) urea;
1-[4 (Piporidine-1-carbonyl)-pyridin-2-yl]-3 (2-pyridin-4-yl-thiazol-4-yl)-urca;
N.N. Diethyl 2 [3 (2 pyridin 4 yl thiazol 4 yl) ureido] isonicotinamide;
1 (4 Diethylaminomethyl-pyridin-2 yl)-3 (2 pyridin-4-yl-thiazol-4-yl) urca;
1-[6-(2,6-Dimethyl-piperidin-1-ylmethyl)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
1-[6-(1-Piperidin-1-yl-ethyl)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
2-(\fo [3-(2-Pyridin 4-yl-thiazol-4-yl)-ureido]-pyridin-2-ylamino\-methyl)-piperidino-1-
     carboxylic acid-tert-butyl-ester;
1-16-[(Piperidin 2 ylmethyl) amino] pyridin 2 yl\-3 (2 pyridin 4 yl thiazol 4 yl) urea;
(S) 1 [6 (3 Methyl-piperidin-1-ylmethyl) pyridin-2 yl] 3 (2 pyridin 4 yl thiazel 4-yl) urea;
(R) 1-[6-(3-Methyl-piperidin-1-ylmethyl)-pyridin-2-yl] 3 (2-pyridin-4-yl-thiazel-4-yl)-urca;
1-(2-Chloro-thiazol 4-yl) 3-(6-piperidin-1-ylmethyl-pyridin-2-yl)-urea;
1-(2-Bromo-thiazol-4-yl)-3-[6-(2-piperidin-4-yl-ethoxy)-pyridin-2-yl]-urea;
1-(2-Chloro-thiazol-4-yl)-3-[6-(2-piperidin-4-yl-ethoxy)-pyridin-2-yl]-urea;
1-[6-(Azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-bromo-thiazol-4-yl)-urea;
1-[6-(Azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-chloro-thiazol-4-yl)-urea;
1-(2-Bromo-thiazol-4-yl)-3-[6-(piperidin-4-ylmethoxy)-pyridin-2-yl]-urea;
1-(2-Chloro-thiazol-4-yl)-3-[6-(piperidin-4-ylmethoxy)-pyridin-2-yl]-urea;
3-(4-{3-[6-(1-Methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-ureido}-thiazol-2-yl)-
     benzenesulfonamide;
tert-Butyl 3-{6-[3-(2-pyridin-4-yl-thiazol-4-yl)-ureido]-pyridin-2-yloxymethyl}-pyrrolidine-1-
     carboxylate;
1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(pyrrolidin-3-ylmethoxy)-pyridin-2-yl]-urea;
1-(2-Cyclopropyl-thiazol-4-yl)-3-[6-(2-piperidin-4-yl-ethoxy)-pyridin-2-yl]-urea;
Isopropyl (6-[3-(2-pyridin 4-yl-thiazol 4-yl) urcido] pyridin 2-ylmethyl) carbamic acid-tert-
     butyl ester;
1-[6 (Isopropylamine methyl) pyridin 2-yl] 3 (2 pyridin 4 yl-thiazol 4-yl) urca;
Isopropyl (6-[3-(2-phenyl-thiazel 4-yl) ureido] pyridin 2-ylmethyl) carbamic acid tert-butyl
     ester;
1-[6-(Isopropylamino-methyl)-pyridin 2-yl]-3-(2-phonyl-thiazol-4-yl)-urea;
1-(2-Bromo-thiazol-4-yl)-3-[6-(1-methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-urea;
1-(2-Chloro-thiazol-4-yl)-3-[6-(1-methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-urea;
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1-(2-phenylthiazol-4-yl)-3-(6-p-pyrrolidin-1-ylmethylphenoxypyridin-2-yl)urea;
1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(tetrahydro-furan-3-yloxy)-pyridin-2-yl]-urea;
1-[2 (1H-Indazel-5-yl) thiazel 4-yl] 3-(6-piperidin-1-ylmethyl-pyridin-2-yl)-urea;
1-(1' Methyl-1',2',3',6' tetrahydro [2,4']bipyridinyl-6-yl)-3-(2-pyridin-4-yl-thiazol-4-yl) urea;
1-(2-Bromo thizol 4-yl) 3-(1'-mothyl-1',2',3',6'-tetrahydro-[2,4']bipyridinyl-6-yl) urou;
1-(1' Methyl-1',2',3',6' tetrahydro-2[2,4]bipyridinyl-6-yl) 3-(2-phonyl-thiazol-4-yl) urea;
1-[6 (3-Hydroxy-propylamino) pyridin-2-yl]-3 (2-pyridin-4-yl-thizel-4-yl) urea;
1-(2-Bromo-thiazol-4-yl)-3-[6(3-hydroxy-propylamino)-pyridin-2-yl]-urea;
1 (1' Mothyl 1', 2', 3', 4', 5', 6' hoxahydro [2,4']bipydrinyl 6 - yl) 3 (2 pyridin 4 yl thiazol 4 yl)
     <del>urca:</del>
1-(1'-Methyl-1',2',3',4',5',6'-hexahydro-[2,4']bipyridinyl-6-yl)-3-(2-phenyl-thiazol-4-yl)-urea;
6-[3-(2-Pyridin 4-yl-thizol-4-yl)-urcide] 3',6' dihydro 2'H-[2,4]bipyridinyl-1' carboxylio acid
     tort-butylester:
1-(2 Pyridin 4 vl thiazol 4 vl) 3-(1',2',3',6' tetrahydro [2,4']bipyridinyl 6-yl) urea;
1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(tetrahydro-furan-3-ylmethoxy)-pyridin-2-yl]-urea;
2-[6-[3-(2-Pyridin-4-yl-thiazol-4-yl)-ureido]-pyridin-2-yloxymethyl]-pyrrolidine-1-carboxylic
     acid tert-butyl ester;
1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-urea;
6-[3-(2-Pyridin 4-yl-thiazol-4-yl)-ureido]-pyridine-2-carbothioic acid-diethylamide;
1-(2-Brome-thiazel-4-yl)-3-[6-(3-methyl-piperidin-1-ylmethyl)-pyridin-2-yl]-urea;
1-(2-Chloro-thiazol-4-yl)-3-[6-(3-methyl-piperidin-1-ylmethyl)-pyridin-2-yl]-urea;
1-[2-(2-Methoxy-pyridin-4-yl)-thiazol-4-yl]-3-[6-(1-methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-
     yl]-urea;
1-[6-(2-Dimethylamino-ethoxy)-pyridin-2-yl]-3-[2-(2-methoxy-pyridin-4-yl)-thiazol-4-yl]-urea;
1-[6-(1-Methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
1-(2-phenylthiazol 4-yl) 3-(6-pyrrolidin-1-ylmethyl-pyridin-2-yl)uroa;
1-(6-Diethylaminomethylpyridin-2-yl)-3-(2-phenylthiazol-4-yl)ureat
(S)-1-[6-(1-Methylpyrrolidin-2-ylmethoxy)pyridin-2-yl]-3-(2-phenylthiazol-4-yl)urea;
1-[6-(2-Piperidin-4-yl-ethoxy)pyridin-2-yl]-3-[2-phenylthiazol-4-yl]urea;
1-[6-(4-Ethylpiperazin-1-yl)-pyridin-2-yl]-3-(2-phenylthiazel-4-yl)urea;
1-(2-phonylthiazol 4-yl)-3-[6-(4-pyrimidin 2-yl-piperazin-1-yl)pyridin 2-yl-proa;
Diethyl 6 [3 (2 phonylthiazol 4 yl)ureide] pyridine 2 carboxamide;
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1-(2-Pyridin-4-yl-thiazol-4-yl)-3-(6-p-pyrrolidin-1-ylmethylphenoxypyridin-2-yl)urea;

1-(2-Bromothiazol-4-yl)-3-(6-p-pyrrolidin-1-ylmethylphenoxypyridin-2-yl)urea;

1-[6-(Piperidin-4-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;

1-[6-(1-Methyl-piperidin-4-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;

1-[6-(1-Methyl-piperidin-4-yloxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;

1-[6-(1-Methyl-azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;

1-[6-(Azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;

1-[6-(1-Methyl-azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;

1-(2-Phenyl-thiazol-4-yl)-3-[6-(piperidin-4-ylmethoxy)-pyridin-2-yl]-urea;

1-[6-(1-Methyl-piperidin-4-ylmethoxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;

1-[6-(1-Methyl-piperidin-4-yloxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;

1-[6-(2-Piperidin-4-yl-ethoxy)-pyridin-2-yl]-3-(2-thiophen-2-yl-thiazol-4-yl)-urea; and

1-[6-(1-Methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-3-[2-(thiophene-2-sulfonylmethyl)-thiazol-4-yl]-urea;

1-[2 (2 Methoxy pyridin 4 yl) thiazol 4 yl] 3 (6 piperdin 1-ylmethyl pyridin 2 yl) urea; and [2 (2 Chloro pyridin 4 yl) thiazol 4 yl] 3 (6 piperdin 1-ylmethyl pyridin 2 yl) urea; and pharmaceutically acceptable salts thereof.

Claim 13 (canceled).

Claim 14 (currently amended): A method of inhibiting cell proliferation which comprises administering an effective amount of a compound of Formula VI

$$\mathbb{R}^{17} \xrightarrow{\begin{array}{c} N \\ 3 \end{array}} \xrightarrow{\begin{array}{c} 4 \\ 1 \end{array}} \xrightarrow{\begin{array}{c} 5 \\ N \end{array}} \xrightarrow{\begin{array}{c} K \\ 15 \end{array}} \mathbb{VI}$$

wherein R¹⁵ is one or more substituents selected from H, optionally substituted heterocyclyl, phonyl, C₁-C₂-alkyl, C₁-C₂-halcalkyl, C₁-C₄-hydroxyalkyl, amino, C₁-C₄-azidoalkyl, C₄-C₄-azidoalkyl, C₄-C₄-C₄-azidoalkyl, C₄-C₄-azidoalkyl

eyanoalkyl, G₄-G₄-aminoalkyl, halo, hydroxy, (optionally substituted heterocyclyl) G₄-G₄-alkyl, optionally substituted phenoxy-G₄-G₂-alkyl, G₄-G₄-alkexy-G₄-G₄-alkyl, G₄-G₄-alkyl, optionally substituted heterocyclyloxy, optionally substituted heterocyclyl-C₁-C₄-alkoxy, C₁-C₄-alkylamino-C₁-C₄-alkoxy, and optionally substituted phenoxy, G₄-G₄-alkoxyearbonyl, 5-6-membered heterocyclyl-G₄-G₄-alkylaminocarbonyl, 5-6-membered N-containing heterocyclylearbonyl, G₄-G₄-alkylaminocarbonyl, G₄-G₄-alkylaminothiocarbonyl, G₄-G₄-alkylaminocarbonyl, 5-6-membered N-containing heterocyclyl-G₄-G₄-alkylaminocarbonyl, aminocarbonyl, 5-6-membered N-containing heterocyclyl-G₄-G₄-alkylamino, G₄-G₄-alkylamino-G₄

wherein R¹⁶ is selected from H, heterocyclylcarbonyl, alkylaminocarbonyl, and alkylaminomethyl, and heterocyclylmethyl; and

wherein R¹⁷ is selected from halo, C₁-C₆-alkyl, cycloalkylalkynyl, cycloalkyl, optionally substituted indelyl, optionally substituted indezelyl, optionally substituted phenexy, optionally substituted heteroarylsulfonyl-G₄-G₄-alkyl thienylsulfonyl-C₁-C₄-alkyl, unsubstituted 5 membered oxygen or sulfur containing heteroaryl, thienyl, unsubstituted 6 membered nitrogen containing heterocyclyl, phenyl optionally substituted with one or two substituents selected

from halo, C₄-C₄-alkylamino, amino, nitro, C₄-C₄-alkoxy, G₄-G₂-haloalkyl, hydroxy, C₄-C₄-alkylthio, C₄-C₄-alkylcarbonylamino, (optionally substituted phenyl)sulfonylamino, eyano, C₄-C₂-haloalkoxy, 5-or-6-membered N-containing heterocyclyl, aminosulfonyl, (6-membered N-containing heterocyclyl)sulfonyl, C₄-C₂-haloalkylcarbonylaminosulfonyl and (optionally substituted phenyl)aminosulfonyl, and 6-membered nitrogen-containing heterocyclyl optionally substituted with one-or-more substituents independently selected from pyridyl, phenyl,

G₁-G₂-alkyl, G₁-G₂-halealkyl, C₁-C₂ alkoxy, amine, hale, piperidinyl, morphelinyl, G₁-G₂-alkylaminethiocarbonyl, N,N-di-G₁-G₂-alkylamine-G₁-G₄-alkylenyl, N-G₁-G₂-alkylamine-G₁-G₄-alkylenyl, morphelinyl-G₁-G₄-alkylenylaminecarbonyl, aminecarbonyl, G₁-G₂-halealkylearbonylamine, morphelinyl-G₁-G₄-alkylenylamine, N,N-di-G₁-G₂-alkylamine and N,N-di-G₁-G₂-alkylamine-G₁-G₄-alkylenylamine;

and pharmaceutically acceptable derivatives salts thereof; provided only one of R¹⁵ and R¹⁶ is H.

Claims 15-66 (Canceled).

Claim 67 (currently amended): The method of Claim 49 111, wherein the compound is and pharmaceutically acceptable salts thereof selected from:

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1-pyridin 2 yl-3 (2 pyridin 4 ylthiazol 4 yl)urca;
1-(6-ethylpyridin-2-yl) 3-(2-pyridin-4-ylthiazol-4-yl)urca;
1-(2-pyridin 4-yl-thiazol-4-yl)-3-(3,4,5,6-tetrahydro-2H-[1,2']bipyridinyl-6'-yl)urca;
1-(6-(diethylaminomethyl)pyridin 2-yl)-3-(2-pyridin 4-ylthiazol 4-yl)urea;
1-[6-(4-methylpiperazin-1-yl)pyridin-2-yl]-3-(2-pyridin-4-ylthiazol-4-yl)urea;
1-[6 (piperidin-1-ylmethyl)pyridin-2-yl]-3-[2-(pyridin-4-yl)thiazol-4-yl]urea;
1-(6-phenoxy-pyridin-2-yl)-3-(2-pyridin-4-yl-thiazol-4-yl)urea;
1-12-(2-ethoxy pyridin 4-yl) thiazol 4-yll-3 (6-ethyl pyridin 2-yl) urea;
1-(6-diethylaminomethyl-pyridin-2-yl)-3 (2-pyridin-3-yl-thiazol-4-yl)-urea;
1-[2-(2-methoxy-pyridin-4-yl)-thiazol-4-yl]-3-(6-morpholin-4-ylmethyl-pyridin-2-yl)-urca;
1-(2-pyridin-4-yl-thiazol-4-yl)-3-(6-pyrrolidin-1-ylmethyl-pyridin-2-yl)-urea;
1-(2-phenylthiazol 4 yl) 3 (6-piperidin-1-ylmethyl-pyridin-2-yl)urea;
1-[6-(1-methylpyrrolidin-2-ylmethoxy)pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)urea; and
1-[2-(4-aminophonyl)thiazol-4-yl]-3-(6-piperidin-1-ylmethyl-pyridin-2-yl)urca; and
1-{6-[4-(2-aminoethyl)phenoxy]pyridin-2-yl}-3-(2-pyridin-4-yl-thiazol-4-yl)urea, and
pharmaceutically acceptable salts thereof.
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Claims 68-110 (canceled)

Claim 111 (currently amended): A method of <u>treating cancer</u> inhibiting cell-preliferation which comprises administering an effective amount of a compound of Formula VI

wherein R¹⁵ is one or more substituents selected from H, optionally substituted heterocyclyl, phenyl, C₁-C₂-alkyl, C₄-C₄-halealkyl, C₄-C₄-hydroxy, (optionally substituted heterocyclyl) C₄-C₄-eyanealkyl, C₄-C₄-aminoalkyl, hale, hydroxy, (optionally substituted heterocyclyl) C₄-C₄-alkyl, optionally substituted phenoxy-C₄-C₂-alkyl, C₄-C₄-alkoxy-C₄-C₄-alkyl, C₄-C₄-alkyl, optionally substituted heterocyclyloxy, optionally substituted heterocyclyl-C₁-C₄-alkoxy, C₁-C₄-alkylamino-C₁-C₄-alkoxy, and optionally substituted phenoxy, C₄-C₄-alkoxycarbonyl, 5 6-membered heterocyclyl-C₄-C₄-alkylaminocarbonyl, 5 6-membered N-containing heterocyclylearbonyl, C₄-C₄-alkylaminocarbonyl, aminocarbonyl, 5 6-membered N-containing heterocyclyl-sulfonyl-C₄-C₄-alkylaminocarbonyl, aminocarbonyl, 5 6-membered N-containing heterocyclyl-sulfonyl-C₄-C₄-alkylamino-C₄-C₄-alkyl

wherein R¹⁶ is selected from H, heterocyclylearbonyl, alkylaminocarbonyl, and alkylaminomethyl, and heterocyclylmethyl; and

wherein R¹⁷ is selected from halo, C₁-C₆-alkyl, cycloalkylalkynyl, cycloalkyl, optionally substituted indazolyl, optionally substituted phonoxy; optionally substituted heteroarylsulfonyl-C₄-G₄-alkyl, thienylsulfonyl-C₁-C₄-alkyl, unsubstituted 5-membered-oxygen or sulfur containing heteroaryl, thienyl, unsubstituted 6-membered nitrogen containing heterocyclyl, phenyl optionally substituted with one or two substituents selected

from halo, C₄-C₄-alkylamino, amino, nitro, C₄-C₄-alkoxy, C₄-C₂-haloalkyl, hydroxy, C₄-C₄-alkylthio, C₄-C₄-alkylcarbonylamino, (optionally substituted phonyl)sulfonylamino, eyano, C₄-C₂-haloalkoxy, 5- or 6-membered N-containing

heterocyclyl, aminosulfonyl, (6-membered-N-containing-heterocyclyl)sulfonyl, C₁-C₂-haloalkylearbonylaminosulfonyl and (optionally substituted-phonyl)aminosulfonyl, and 6-membered nitrogen-containing heterocyclyl optionally substituted with one or more substituents independently selected from pyridyl, phonyl,

G₁-G₂-alkylpiperazinyl, G₁-G₂-alkylaminethiocarbonyl, N,N-di-G₁-G₂-alkylamine-G₁-G₄-alkylenyl, N-G₄-G₂-alkylamine-G₄-alkylenyl, morpholinyl-G₄-G₄-alkylenyl, aminocarbonyl, G₄-G₂-haloalkylearbonylamine, morpholinyl-G₄-G₄-alkylenylamine, N,N-di-G₄-G₂-haloalkylearbonylamine, and N,N-di-G₄-G₂-alkylamine and N,N-di-G₄-G₂-alkylamine G₄-G₄-alkylenylamine;

and pharmaceutically acceptable derivatives salts thereof; provided only one of R¹⁵ and R¹⁶ is H.

Claim 112 (currently amended): The method of Claim 111, wherein R¹⁵ is selected from H, optionally substituted pyrrolidinyl, optionally substituted piperazinyl, optionally substituted piperidinyl, merehelinyl, 1.2.3.6 tetrahydro-pyridinyl, (optionally substituted pyrrolidinyl) - Gu-C2-alkyl. (optionally substituted piperidinyl) C1-C2-alkyl, (optionally substituted piperazinyl) C1-C2-alkyl, morpholinyl-C1-C2-alkyl, C1-C4-alkylamino-C1-C4-alkyl, C1-C4-hydroxyalkylamino, (optionally substituted pyrrolidinyl) G. G. alkylamino; (optionally substituted piperidinyl) -G. G2-alkylamino, (optionally substituted-piperazinyl) G1 G2-alkylamino, morpholinyl G1-G2alkylamino, optionally substituted pyrrolidinyl-C₁-C₄-alkoxy, optionally substituted azetidinyl-C₁-C₄-alkoxy, tetrahydrofuryl-C₁-C₄-alkoxy, optionally substituted piperidinyl-C₁-C₄-alkoxy, C₁-C₄-alkylamino-C₁-C₄-alkoxy, tetrahydrofuryloxy, optionally substituted piperidinyloxy, and optionally substituted phenoxy. G. G. alkylaminocarbonyl and G. G. alkylaminothiocarbonyl; wherein R¹⁶ is selected from H, 5-6-membered nitrogen-containing heterocyclylearbonyl, C₁-C₄alkylaminocarbonyl, and C₁-C₄-alkylaminomethyl, and 5-6-membered-nitrogen-containing heteroeyelylmethyl; and wherein R¹⁷ is selected from halo, C₁-C₂-alkyl, optionally substituted 5-6-membered heteroarylsulfonyl-C1-C2-alkyl, thienylsulfonyl-C1-C2-alkyl, optionally substituted phenoxy; and C₃-C₆-cycloalkyl-C₂-C₄-alkynyl; and pharmaceutically acceptable derivatives thereof.

Claim 113 (currently amended): The method of Claim 112, wherein R¹⁵ is selected from H. tetrahydro-furanyloxy, 1-methylpyrrolidin-2-ylmethoxy, 2-pyrrolidinylmethoxy, 3pyrrolidinylmethoxy, 1-Boc-pyrrolidin-2-ylmethoxy, 4-piperidinylmethoxy, 1-Boc-piperidin-4ylmethoxy, 1-Boc-piperidin-4-ylethoxy, piperidin-4-ylethoxy, 1-methyl-piperidin-4-ylmethoxy, 1-Boc-azetidin-3-ylmethoxy, 1-methyl-azetidin-3-ylmethoxy, 3-azetidinylmethoxy, 1-methylpiperidin-4-yloxy, phenyloxy, phenoxy, 4-(pyrrolidin-1-ylmethyl)phenoxy, and dimethylaminoethoxy, 1-piperidinylmethyl, 1-(piperidin-1-yl)ethyl, 3-methylpiperidin-1vimethyl, 1 pyrrolidinylmethyl, 2,2,6,6 tetramethylpiperidin 1 ylmethyl, 2,6 dimethylpiperidin 1-ylmethyl, dimethylaminemethyl, diethylaminemethyl, diethylaminethiocarbenyl, diethylaminocarbonyl, N. Boe N. isopropylaminomethyl, isopropylaminomethyl, 2 thionylculfonylmothyl, hydroxypropylamino, 4 othyl-piperidin 1-yl, 4-(2-pyridyl)piperidin 1-yl, 1-methylpiperidin 4 vl. 4 (2-pyrazinyl)piperidin 1-yl, 1-methyl-1,2,3,6-tetrahydro-pyridin 4-yl, 1,2,3,6-tetrahydro-pyridin 4-yl, and 1-Boo-1,2,3,6-tetrahydro-pyridin 4-yl; wherein R¹⁶ is selected from H, 1-piperidinylearbonyl, diethylaminocarbonyl, and diethylaminomethyl, 1-piperidinylmethyl: and wherein R¹⁷ is selected from chloro, bromo, methyl and cyclopropylethynyl and pharmaceutically acceptable derivatives thereof.

Claim 114 (currently amended): The method of Claim 113, wherein R¹⁷ is chloro or bromo; and pharmaccutically acceptable derivatives thereof.

Claim 115 (currently amended): The method of Claim 111, wherein R¹⁵ is selected from H, optionally substituted pyrrolidinyl, optionally substituted piperazinyl, optionally substituted piperazinyl, (optionally substituted pyrrolidinyl) C₁-G₂-alkyl, (optionally substituted piperazinyl) C₁-G₂-alkyl, (optionally substituted piperazinyl) C₂-G₃-alkyl, morpholinyl C₄-G₂-alkyl, (optionally substituted piperazinyl) C₄-G₅-alkylamino, (optionally substituted piperazinyl) C₄-G₅-alkylamino, morpholinyl C₄-C₅-alkylamino, C₄-C₄-alkylamino-C₄-G₄-alkyl₂-C₄-hydroxyalkylamino, optionally substituted pyrrolidinyl-C₁-C₄-alkoxy, optionally substituted azetidinyl-C₁-C₄-alkoxy, tetrahydrofuryl-C₁-C₄-alkoxy, optionally substituted piperidinyl-C₁-C₄-alkoxy, optionally substituted piperidinyloxy, optionally substituted piperidinyloxy

heterocyclylearbonyl, C₁-C₄-alkylaminocarbonyl, and C₁-C₄-alkylaminomethyl, and 5-6-membered nitrogen containing heterocyclylmethyl; and wherein R¹⁷ is selected from C₃-C₆-cycloalkyl and phenyl-optionally substituted with one or two substituents selected from halo; C₄-C₄-alkylamino; amino; nitro; C₄-C₄-alkoxy; C₄-C₂-haloalkyl, hydroxy; C₄-C₄-alkylthio, C₄-C₄-alkylcarbonylamino, (optionally substituted phenyl)sulfonylamino; cyano; C₄-C₂-haloalkoxy, 5- or 6-membered N-containing heterocyclyl, aminosulfonyl, (6-membered N-containing heterocyclyl)sulfonyl, C₄-C₂-haloalkylcarbonylaminosulfonyl and (optionally substituted phenyl)aminosulfonyl;

and pharmaceutically acceptable derivatives thereof.

Claim 116 (currently amended): The method of Claim 115, wherein R¹⁵ is selected from H, tetrahydro-furanyloxy, 1-methylpyrrolidin-2-ylmethoxy, 2-pyrrolidinylmethoxy, 3pyrrolidinylmethoxy, 1-Boc-pyrrolidin-2-ylmethoxy, 4-piperidinylmethoxy, 1-Boc-piperidin-4vlmethoxy, 1-Boc-piperidin-4-vlethoxy, piperidin-4-vlethoxy, 1-methyl-piperidin-4-vlmethoxy, 1-Boc-azetidin-3-ylmethoxy, 1-methyl-azetidin-3-ylmethoxy, 3-azetidinylmethoxy, 1-methylpiperidin-4-yloxy, phenyloxy, phenoxy, 4-(pyrrolidin-1-ylmethyl)phenoxy, and dimethylaminoethoxy, 1-piperidinylmethyl, 1-(piperidin-1-yl)cthyl, 3-methylpiperidin-1ylmothyl, 1-pyrrolidinylmothyl, 2,2,6,6-tetramethylpiperidin-1-ylmothyl, 2,6-dimethylpiperidin-1-vlmethyl.-dimethylaminomethyl.-diethylaminomethyl, diethylaminothiocarbonyl, diethylaminocarbonyl. N. Boe. N. isopropylaminomethyl. isopropylaminomethyl. 2thicnylsulfonylmethyl, hydroxypropylamino, 4-ethyl-piperidin-1-yl, 4-(2-pyridyl)piperidin-1-yl, 1-mothylpiperidin 1 yl, 1 (2 pyrazinyl)piperidin 1 yl, 1-methyl-1,2,3,6-tetrahydro-pyridin 4-yl, 1,2,3,6 tetrahydro-pyridin 4 yl, and 1-Box 1,2,3,6 tetrahydro-pyridin 4 yl; wherein R¹⁶ is selected from H. 1-piperidinylearbonyl, diethylaminocarbonyl, and diethylaminomethyl, 1piperidinylmethyl; and wherein R¹⁷ is selected from cyclopropyl and phenyl eptionally substituted with aminosulfonyl: and pharmaccutically acceptable derivatives thereof.

Claim 117 (currently amended): The method of Claim 116, wherein R¹⁷ is unsubstituted phenyls, and pharmaceutically acceptable derivatives thereof.

Claim 118 (currently amended): The method of Claim 111, wherein R¹⁵ is selected from H, optionally substituted pyrrolidinyl, optionally substituted piperazinyl, optionally substituted

piperidinyl, morpholinyl, 1,2,3,6 totrahydro-pyridinyl, (optionally substituted pyrrolidinyl) Ci-C2-alkyl, (optionally substituted piperidinyl) C4-C2-alkyl, (optionally substituted piperazinyl) C4-Ga-alkyl. morpholinyl-Ga-alkyl. (optionally substituted pyrrolidinyl) Ga-Ga-alkylamino, (optionally substituted piperidinyl)-C₁-C₂-alkylamino, (optionally substituted piperazinyl)-C₁-C₂alkylamino, morpholinyl-G₁-G₂-alkylamino, G₁-G₄-alkylamino-G₁-G₄-alkyl, G₁-G₄hydroxyalkylamino, optionally substituted pyrrolidinyl-C₁-C₄-alkoxy, optionally substituted azetidinyl-C₁-C₄-alkoxy, tetrahydrofuryl-C₁-C₄-alkoxy, optionally substituted piperidinyl-C₁-C₄alkoxy, C₁-C₄-alkylamino-C₁-C₄-alkoxy, tetrahydrofuryloxy, optionally substituted piperidinyloxy, and optionally substituted phenoxy. G₁-G₂-alkylaminocarbonyl and G₁-G₂-G₄alkylaminethiocarbonyl: wherein R¹⁶ is selected from H, 5-6-membered nitrogen containing heterocyclylearbonyl, C₁-C₄-alkylaminocarbonyl, and C₁-C₄-alkylaminomethyl, and 5-6membered nitrogen-containing heterocyclylmethyl; and wherein R¹⁷ is selected from optionally substituted indexelyl, optionally substituted indelyl, unsubstituted-5 membered exygen or sulfur containing heteroaryl, unsubstituted thienyl, unsubstituted 6-membered nitrogen-containing heterocyclyl, and 6-membered nitrogen-containing heterocyclyl substituted with ene or more substituents independently selected from pyridyl, phonyl,

G₄-G₄-alkyl, G₄-G₂-haloalkyl, C₁-C₂ alkoxy, amino, halo, piperidinyl, morpholinyl, G₄-G₂ alkylamino-G₄-G₄-alkylamino-G₄-G₄-alkylamino-G₄-G₄-alkylamino-G₄-G₄-alkylamino-G₄-G₄-alkylamino-G₄-G₄-alkylamino-G₄-G₄-haloalkylcarbonylamino, morpholinyl-G₄-G₄-alkylamino, N;N-di-G₄-G₂-alkylamino-G₄-G₄-alkylamino-G₄-G₄-alkylamino;

and pharmaccutically acceptable derivatives thereof.

Claim 119 (currently amended): The method of Claim 118, wherein R¹⁵ is selected from H, tetrahydro-furanyloxy, 1-methylpyrrolidin-2-ylmethoxy, 2-pyrrolidinylmethoxy, 3-pyrrolidinylmethoxy, 1-Boc-pyrrolidin-2-ylmethoxy, 4-piperidinylmethoxy, 1-Boc-piperidin-4-ylmethoxy, 1-methyl-piperidin-4-ylmethoxy, 1-methyl-piperidin-4-ylmethoxy, 1-methyl-piperidin-4-ylmethoxy, 1-methyl-piperidin-4-yloxy, phenyloxy, 4-(pyrrolidin-1-ylmethyl)phenoxy, and dimethylaminoethoxy, 1-pyrrolidinylmethyl, 1-pyrrolidinylmethyl, 2,2,6,6-tetramethylpiperidin-1-ylmethyl, 2,6-dimethylpiperidin-1-ylmethyl,

dimethylaminomethyl, diethylaminomethyl, diethylaminothiocarbonyl, diethylaminocarbonyl, N Boe N isopropylaminomethyl, isopropylaminomethyl, 2-thionylsulfonylmethyl, hydroxypropylamino, 4-ethyl-piperidin 1-yl, 4-(2-pyridyl)piperidin 1-yl, 1-methyl-piperidin 4-yl, 4-(2-pyrazinyl)piperidin 1-yl, 1-methyl-1,2,3,6-tetrahydro-pyridin 4-yl, 1,2,3,6-tetrahydro-pyridin 4-yl, 1,2,3,6-tetrahydro-pyridin 4-yl, and 1-Boe 1,2,3,6-tetrahydro-pyridin 4-yl; wherein R¹⁶ is selected from H, 1-piperidinylearbonyl, diethylaminocarbonyl, and diethylaminomethyl, 1-piperidinylmethyl; and wherein R¹⁷ is selected from 5-indazelyl, 1-Boe indol-5-yl, unsubstituted thienyl, 5-tert-butyloxazel-2-yl and 4-pyridyl substituted with one-or-more substituents independently selected from methoxy-and-chloro; and pharmaceutically acceptable derivatives thereof.

and pharmaceutically acceptable derivatives incredi.

Claim 120 (currently amended): The method of Claim 119 111, wherein R¹⁷ is 4-pyridyl; and pharmaceutically acceptable derivatives thereof.

Claim 121 (currently amended): The method of Claim 111, wherein the compound is and pharmaceutically acceptable derivatives thereof selected from:

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1-[6-(3-Mothyl-piperidin-1-ylmethyl) pyridin 2-yl] 3-(2-phenyl-thiazol-4-yl) urea;
1-[4-(Piperidine-1-earbonyl) pyridin 2-yl] 3-(2-pyridin-1-yl-thiazol-4-yl) urea;
1-(2-Chloro-thiazol-4-yl) 3-[4-(piperidine-1-earbonyl) pyridin-2-yl] urea;
N,N-Diethyl-2-[3-(2-pyridin-4-yl-thiazol-4-yl) ureido] isonicotinamide;
N,N-Diethyl-2-[3-(2-phenyl-thiazol-4-yl) ureido] isonicotinamide;
2-[3-(2-Bromo-thiazol-4-yl) ureido] N,N-diethyl-isonicotinamide;
1-(4-Diethylaminomethyl-pyridin-2-yl) 3-(2-pyridin-4-yl-thiazol-4-yl) urea;
1-[6-(2,6-Dimethyl-piperidin-1-ylmethyl) pyridin-2-yl] 3-(2-pyridin-4-yl-thiazol-4-yl) urea;
1-[6-(1-Piperidin-1-yl-ethyl) pyridin-2-yl] 3-(2-pyridin-4-yl-thiazol-4-yl) urea;
2-({6-[3-(2-Pyridin-4-yl-thiazol-4-yl) ureido] pyridin-2-yl] 3-(2-pyridin-4-yl-thiazol-4-yl) urea;
(S) 1-[6-(3-Mothyl-piperidin-1-ylmethyl) pyridin-2-yl] 3-(2-pyridin-4-yl-thiazol-4-yl) urea;
(R) 1-[6-(3-Mothyl-piperidin-1-ylmethyl) pyridin-2-yl] 3-(2-pyridin-4-yl-thiazol-4-yl) urea;
(R) 1-[6-(3-Mothyl-piperidin-1-ylmethyl) pyridin-2-yl] 3-(2-pyridin-4-yl-thiazol-4-yl) urea;
(R) 1-[6-(3-Mothyl-piperidin-1-ylmethyl) pyridin-2-yl] 3-(2-pyridin-4-yl-thiazol-4-yl) urea;
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1-(2-Bromo-thiazol-4-yl)-3-[6-(2-piperidin-4-yl-ethoxy)-pyridin-2-yl]-urea;
1-(2-Chloro-thiazol-4-yl)-3-[6-(2-piperidin-4-yl-ethoxy)-pyridin-2-yl]-urea;
1-[6-(Azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-bromo-thiazol-4-yl)-urea;
1-[6-(Azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-chloro-thiazol-4-yl)-urea;
1-(2-Bromo-thiazol-4-yl)-3-[6-(piperidin-4-ylmethoxy)-pyridin-2-yl]-urea;
1-(2-Chloro-thiazol-4-yl)-3-[6-(piperidin-4-ylmethoxy)-pyridin-2-yl]-urea;
tert-Butyl 3-{6-[3-(2-pyridin-4-yl-thiazol-4-yl)-ureido]-pyridin-2-yloxymethyl}-pyrrolidine-1-
     carboxylate;
1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(pyrrolidin-3-ylmethoxy)-pyridin-2-yl]-urea;
1-(2-Cyclopropyl-thiazol-4-yl)-3-[6-(2-piperidin-4-yl-ethoxy)-pyridin-2-yl]-urea;
1 [6 (Isopropylamino methyl) pyridin 2 yl] 3 (2 pyridin 4 yl thiazol 4 yl) urea;
1-[6 (Isopropylamino methyl) pyridin 2 yl] 3 (2 phonyl thiazol 4 yl) urea;
1-(2 Bromo-thiazol 4-yl)-3-[6 (isopropylamino-methyl) pyridin-2-yl]-urea;
1-(2-Bromo-thiazol-4-yl)-3-[6-(1-methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-urea;
1-(2-Chloro-thiazol-4-yl)-3-[6-(1-methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-urea;
1-(2-phenylthiazol-4-yl)-3-(6-p-pyrrolidin-1-ylmethylphenoxypyridin-2-yl)urea;
1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(tetrahydro-furan-3-yloxy)-pyridin-2-yl]-urea;
1-[2 (1H-Indazol-5 yl) thiazol-4 yl] 3 (6 piperidin-1-ylmethyl-pyridin-2-yl) urea;
1-(1' Methyl 1',2',3',6' tetrahydro [2,4']bipyridinyl 6 yl) 3 (2 pyridin 4 yl thiazol 4 yl) urca;
1-(2-Bromo thizol-4-yl)-3-(1'-methyl-1',2',3',6' tetrahydro-[2,4']bipyridinyl-6-yl)-urca;
1-(1'-Methyl-1',2',3',6' tetrahydro-2[2,4]bipyridinyl-6-yl) 3-(2-phenyl-thiazol-4-yl) urea;
1 [6 (3 Hydroxy propylamino) pyridin 2 yl] 3 (2 pyridin 4 yl thizol 4 yl) urea;
1 (2 Brome thiazel 4 yl) 3 [6(3 hydroxy propylamine) pyridin 2 yl] urea;
1 (1' Methyl 1',2',3',4',5',6' hexahydro [2,4']bipydrinyl 6 yl) 3 (2 pyridin 4 yl thiazol 4 yl)
     urca;
1 (1' Mothyl-1',2',3',4',5',6' hexahydro [2,4']bipyridinyl-6 yl)-3 (2 phonyl-thiazol-4
6 [3 (2 Pyridin 4 yl thizol 4 yl) uroido] 3',6' dihydro 2'H [2,4]bipyridinyl-1' carboxylic acid
     tert-butylester;
1 (2 Pyridin 4 vl thiazol 4 vl) 3 (122.3262 tetrahydro [2,42]bipyridinyl 6 vl) urca;
1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(tetrahydro-furan-3-ylmethoxy)-pyridin-2-yl]-urea;
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- 2-[6-[3-(2-Pyridin-4-yl-thiazol-4-yl)-ureido]-pyridin-2-yloxymethyl]-pyrrolidine-1-carboxylic acid tert-butyl ester;
- 1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-urea;
- 6 [3 (2 Pyridin 4 yl thiazol 4 yl) uroido] pyridino 2 carbothioic acid diethylamide;
- 1-(2-Bromo-thiazol 4-yl)-3-[6-(3-methyl-piperidin-1-ylmethyl)-pyridin-2-yl]-urea;
- 1-(2-Chloro-thiazol-4-yl)-3-[6-(3-methyl-piperidin-1-ylmethyl)-pyridin-2-yl]-urea;
- 1 (2 Phenyl-thiazol 4 yl) 3 [4 (piperidine-1 carbonyl) pyridin-2 yl] urea;
- 1-(2-Bromo-thiazol 4-yl)-3-[4-(piperidine-1-carbonyl)-pyridin-2-yl]-urea;
- 1-[2-(2-Methoxy-pyridin-4-yl)-thiazol-4-yl]-3-(6-phenoxy-pyridin-2-yl)-urea;
- 1-[2-(2-Methoxy-pyridin-4-yl)-thiazol-4-yl]-3-[6-(1-methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-urea;
- 1-[6-(2-Dimethylamino-ethoxy)-pyridin-2-yl]-3-[2-(2-methoxy-pyridin-4-yl)-thiazol-4-yl]-urea;
- 1-[6-(1-Methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-(2-phenylthiazol 4-yl)-3 (6-pyrrolidin-1-ylmethyl-pyridin-2-yl)urea;
- 1-(6-Diethylaminemethylpyridin-2-yl)-3-(2-phenylthiazol-4-yl)urca;
- (S)-1-[6-(1-Methylpyrrolidin-2-ylmethoxy)pyridin-2-yl]-3-(2-phenylthiazol-4-yl)urea;
- 1-[6-(2-Piperidin-4-yl-ethoxy)pyridin-2-yl]-3-[2-phenylthiazol-4-yl]urea;
- 1-[6-(4-Ethylpiperazin-1-yl)-pyridin-2-yl]-3-(2-phenylthiazol-4-yl)urea;

Diethyl 6-[3-(2-phenylthiazol-4-yl)ureido]-pyridine-2-carboxamide;

- 1-(2-Pyridin-4-yl-thiazol-4-yl)-3-(6-p-pyrrolidin-1-ylmethylphenoxypyridin-2-yl)urea;
- 1-(2-Bromothiazol-4-yl)-3-(6-p-pyrrolidin-1-ylmethylphenoxypyridin-2-yl)urea;
- 1-[6-(Piperidin-4-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-piperidin-4-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-piperidin-4-yloxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-[6-(Azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;
- 1-(2-Phenyl-thiazol-4-yl)-3-[6-(piperidin-4-ylmethoxy)-pyridin-2-yl]-urea;
- 1-[6-(1-Methyl-piperidin-4-ylmethoxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-piperidin-4-yloxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;
- 1-[6-(2-Piperidin-4-yl-ethoxy)-pyridin-2-yl]-3-(2-thiophen-2-yl-thiazol-4-yl)-urea; and

- 1-[6-(1-Methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-3-[2-(thiophene-2-sulfonylmethyl)-thiazol-4-yl]-urea;
- 1-[2 (2 Methoxy pyridin 4 yl) thiazel 4 yl] 3 (6 piperdin 1 ylmethyl pyridin 2 yl) urea; and [2 (2 Chloro pyridin 4 yl) thiazel 4 yl] 3 (6 piperdin 1 ylmethyl pyridin 2 yl) urea phamaceutucally acceptable salts thereof.

Claim 122 (currently amended): The method of Claim 111, wherein the compound is and pharmaceutically acceptable derivatives thereof selected from:

- 1-f6-(3-Methyl-piperidin-1-ylmethyl)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;
- 1-[4-(Piperidine-1-carbonyl)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazel-4-yl)-urca;
- N.N. Diethyl-2-[3-(2-pyridin-4-yl-thiazol-4-yl) urcido] isonicotinamido;
- 4 (4 Diethylaminomethyl pyridin 2-yl) 3 (2-pyridin 4-yl thiazol 4-yl) urea;
- 1-[6-(2,6-Dimothyl-piperidin-1-ylmethyl) pyridin-2-yl] 3-(2-pyridin-4-yl-thiazol-4-yl) urea;
- 1-[6 (1 Piperidin 1-yl-ethyl) pyridin 2-yl] 3 (2-pyridin 4-yl-thiazol 4-yl) urea;
- 2-({6-[3 (2-Pyridin 4 yl thiazol 4 yl) urcido] pyridin 2 ylamino} methyl) piperidine-1carboxylic acid tert-butyl-ester;
- 1-{6-[(Piperidin-2-ylmethyl)-amine]-pyridin-2-yl}-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- (S) 1 [6 (3 Mothyl piperidin 1 ylmethyl) pyridin 2 yl] 3 (2 pyridin 4 yl thiazol 4 yl) urea;
- (R) 1 [6 (3 Methyl piperidin 1 ylmothyl) pyridin 2 yl] 3 (2 pyridin 4 yl thiazol 4 yl) uroa;
- 1-(2 Chloro thiazol 4-yl) 3-(6-piperidin-1-ylmethyl-pyridin-2-yl) urea;
- 1-(2-Bromo-thiazol-4-yl)-3-[6-(2-piperidin-4-yl-ethoxy)-pyridin-2-yl]-urea;
- 1-(2-Chloro-thiazol-4-yl)-3-[6-(2-piperidin-4-yl-ethoxy)-pyridin-2-yl]-urea;
- 1-[6-(Azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-bromo-thiazol-4-yl)-urea;
- 1-[6-(Azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-chloro-thiazol-4-yl)-urea;
- 1-(2-Bromo-thiazol-4-yl)-3-[6-(piperidin-4-ylmethoxy)-pyridin-2-yl]-urea;
- 1-(2-Chloro-thiazol-4-yl)-3-[6-(piperidin-4-ylmethoxy)-pyridin-2-yl]-urea;
- 3-(4-{3-[6-(1-Methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-ureido}-thiazol-2-yl)-benzenesulfonamide;
- *tert*-Butyl 3-{6-[3-(2-pyridin-4-yl-thiazol-4-yl)-ureido]-pyridin-2-yloxymethyl}-pyrrolidine-1-carboxylate;
- 1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(pyrrolidin-3-ylmethoxy)-pyridin-2-yl]-urea;

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1-(2-Cyclopropyl-thiazol-4-yl)-3-[6-(2-piperidin-4-yl-ethoxy)-pyridin-2-yl]-urea;

Isopropyl-{6 [3 (2 pyridin 4 yl-thiazol 4 yl) ureido] pyridin 2 ylmethyl}-carbamic acid tert-butyl-ester;

1-[6 (Isopropylamino-methyl) pyridin-2-yl] 3 (2 pyridin-4-yl-thiazol 4-yl) urea;

Isopropyl-{6 [3 (2 phenyl-thiazol 4-yl) ureido] pyridin-2 ylmethyl}-carbamic acid tert-butyl
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1-[6 (Isopropylamino methyl) pyridin 2-yl] 3 (2-phonyl-thiazol-4-yl) urea;

ester:

- 1-(2-Bromo-thiazol-4-yl)-3-[6-(1-methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-urea;
- 1-(2-Chloro-thiazol-4-yl)-3-[6-(1-methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-urea;
- 1-(2-phenylthiazol-4-yl)-3-(6-p-pyrrolidin-1-ylmethylphenoxypyridin-2-yl)urea;
- 1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(tetrahydro-furan-3-yloxy)-pyridin-2-yl]-urea;
- 4-[2-(1H-Indazol-5-yl)-thiazol-4-yl]-3-(6-piperidin-1-ylmethyl-pyridin-2-yl)-urea;
- 4 (1'-Mothyl-1',2',3',6'-tetrahydro-[2,4']bipyridinyl-6-yl)-3 (2-pyridin-4-yl-thiazol-4-yl)-urca;
- 1-(2 Bromo-thizol 4-yl) 3-(1'-methyl-1',2',3',6'-tetrahydro-[2,4']bipyridinyl-6-yl) urca;
- 1-(1'-Methyl-1',2',3',6'-tetrahydro-2[2,4]bipyridinyl-6-yl)-3-(2-phenyl-thiazol-4-yl)-urea;
- 1-16 (3 Hydroxy propylamino) pyridin-2 yl-3 (2 pyridin-4 yl-thizol-4 yl) urea;
- 1-(2-Bromo thiazol 4-yl) 3-[6(3-hydroxy-propylamino)-pyridin-2-yl]-urea;
- 1 (1' Methyl 1',2',3',4',5',6' hexaltydro [2,4']bipydrinyl 6 yl) 3 (2 pyridin 4 yl thiazol 4 yl) urca;
- 1 (1' Methyl-1',2',3',4',5',6' hexahydro [2,4']bipyridinyl-6-yl) 3 (2-phonyl-thiazol-4-yl) urea;
- 6 [3 (2 Pyridin 4 yl thizel 4 yl) ureide] 3',6' dihydre 2'H [2,4]bipyridinyl-1' carboxylic acid tert-butylester;
- 1-(2 Pyridin 4 yl-thiazol 4 yl) 3-(1',2',3',6'-tetrahydro [2,4']bipyridinyl-6-yl) urca;
- 1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(tetrahydro-furan-3-ylmethoxy)-pyridin-2-yl]-urea;
- 2-[6-[3-(2-Pyridin-4-yl-thiazol-4-yl)-ureido]-pyridin-2-yloxymethyl]-pyrrolidine-1-carboxylic acid tert-butyl ester;
- 1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-urea;
- 6-[3-(2-Pyridin-4-yl-thiazol-4-yl) uroido]-pyridino-2-carbothioic acid-diothylamide;
- 1 (2 Brome thiazel 4 yl) 3 [6 (3 mothyl piperidin 1 ylmethyl) pyridin 2 yl] urea;
- 1-(2 Chloro-thiazol 4 yl) 3-[6 (3 methyl piperidin 1 ylmethyl) pyridin 2 yl] urea;
- 1-[2-(2-Methoxy-pyridin-4-yl)-thiazol-4-yl]-3-[6-(1-methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-urea;

- 1-[6-(2-Dimethylamino-ethoxy)-pyridin-2-yl]-3-[2-(2-methoxy-pyridin-4-yl)-thiazol-4-yl]-urea;
- 1-[6-(1-Methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1 (2 phonylthiazol 4 yl) 3 (6 pyrrolidin 1 ylmethyl pyridin 2 yl)urea;
- 4 (6 Diethylaminomethylpyridin 2 yl) 3 (2 phenylthiazol 4 yl)urea;
- (S)-1-[6-(1-Methylpyrrolidin-2-ylmethoxy)pyridin-2-yl]-3-(2-phenylthiazol-4-yl)urea;
- 1-[6-(2-Piperidin-4-yl-ethoxy)pyridin-2-yl]-3-[2-phenylthiazol-4-yl]urea;
- 1-[6-(4-Ethylpiperazin-1-yl)-pyridin-2-yl]-3-(2-phonylthiazol-4-yl)urea;
- 1-(2-phenylthiazol 4-yl)-3-[6-(4-pyrimidin-2-yl-piperazin-1-yl)pyridin-2-yl]urea;

Diethyl-6 [3 (2 phenylthiazol-4-yl)urcido]-pyridine-2-carboxamide;

- 1-(2-Pyridin-4-yl-thiazol-4-yl)-3-(6-p-pyrrolidin-1-ylmethylphenoxypyridin-2-yl)urea;
- 1-(2-Bromothiazol-4-yl)-3-(6-p-pyrrolidin-1-ylmethylphenoxypyridin-2-yl)urea;
- 1-[6-(Piperidin-4-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-piperidin-4-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-piperidin-4-yloxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-[6-(Azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;
- 1-(2-Phenyl-thiazol-4-yl)-3-[6-(piperidin-4-ylmethoxy)-pyridin-2-yl]-urea;
- 1-[6-(1-Methyl-piperidin-4-ylmethoxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-piperidin-4-yloxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;
- 1-[6-(2-Piperidin-4-yl-ethoxy)-pyridin-2-yl]-3-(2-thiophen-2-yl-thiazol-4-yl)-urea; and
- 1-[6-(1-Methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-3-[2-(thiophene-2-sulfonylmethyl)-thiazol-4-yl]-urea;
- 1-[2 (2 Methoxy pyridin 1 yl) thiazol 1 yl] 3 (6 piperdin 1 ylmethyl pyridin 2 yl) urea; and [2 (2 Chloro pyridin 1 yl) thiazol 1 yl] 3 (6 piperdin 1 ylmethyl pyridin 2 yl) urea; and pharmaceutically acceptable salts thereof.
- Claim 123 (currently amended): A method of inhibiting a serine/threonine kinase which comprises administering an effective amount of a compound of Formula VI

$$\mathbb{R}^{17} \xrightarrow{\begin{array}{c} 1 \\ 3 \\ 2 \\ 1 \\ 5 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \\ 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow{\begin{array}{c} 1 \\ 1 \end{array}} \mathbb{N} \xrightarrow$$

wherein R¹⁵ is one or more substituents selected from H, optionally substituted heterocyclyl, phenyl, C₄-C₂-alkyl, C₄-C₂-halcalkyl, C₄-C₄-hydroxyalkyl, amino, C₄-C₄-azidealkyl, C₄-C₄-eyanoalkyl, C₄-C₄-aminoalkyl, halo, hydroxy, (optionally substituted heterocyclyl) C₄-C₄-alkyl, optionally substituted phenoxy-C₄-C₂-alkyl, C₄-C₄-alkoxy-C₄-C₄-alkyl, C₄-C₄-alkyl, optionally substituted heterocyclyloxy, optionally substituted heterocyclyl-C₁-C₄-alkoxy, C₁-C₄-alkylamino-C₁-C₄-alkoxy, and optionally substituted phenoxy, C₄-C₄-alkoxyearbonyl, 5-6-membered N-containing heterocyclyl-C₄-C₄-alkylaminocarbonyl, C₄-C₄-alkylaminothiocarbonyl, C₄-C₄-alkylaminocarbonyl, aminocarbonyl, 5-6-membered N-containing heterocyclyl-sulfonyl-C₄-C₄-alkylamino-C₄-C₄

wherein R¹⁶ is selected from H, heterocyclylearbonyl, alkylaminocarbonyl, and alkylaminomethyl, and heterocyclylmethyl; and

wherein R¹⁷ is selected from halo, C₁-C₆-alkyl, cycloalkylalkynyl, cycloalkyl, optionally substituted indazelyl, optionally substituted phenoxy; optionally substituted heteroarylsulfonyl-C₁-G₄-alkyl thienylsulfonyl-C₁-C₄-alkyl, unsubstituted 5-membered oxygen or sulfur-containing heteroaryl; thienyl, unsubstituted 6-membered nitrogen containing heterocyclyl, phenyl optionally substituted with one or two substituents selected

from halo, C₄-C₄-alkylamino, amino, nitro, C₄-C₄-alkoxy, C₄-C₂-haloalkyl, hydroxy, C₄-C₄-alkylthio, C₄-C₄-alkylearbonylamino, (optionally substituted phenyl)sulfonylamino, evano, C₄-C₂-haloalkoxy, 5-or 6-membered N-containing

heterocyclyl, aminosulfonyl, (6-membered N-centaining heterocyclyl)sulfonyl, C₁-C₂-halcalkylearbonylaminosulfonyl and (optionally substituted phenyl)aminosulfonyl, and 6-membered nitrogen-containing heterocyclyl optionally substituted with one or more substituents independently selected from pyridyl, phenyl,

G₁-G₂-alkylpiperazinyl, G₁-G₂-alkylaminothiocarbonyl, N,N-di-G₁-G₂-alkylamino-G₁-G₄-alkylenyl, N-G₄-G₂-alkylamino-G₄-G₄-alkylenyl, morpholinyl-G₄-G₄-alkylenyl, morpholinyl-G₄-G₄-alkylenyl, morpholinyl-G₄-G₄-alkylenylamino-and-N,N-di-G₄-G₂-alkylamino-and-N,N-di-G₄-G₂-alkylamino-and-N,N-di-G₄-G₂-alkylamino-and-N,N-di-G₄-G₂-alkylamino-G₄-G₄-alkylenylamino;

and pharmaceutically acceptable derivatives salts thereof; provided only one of R¹⁵ and R¹⁶ is H.

Claim 124 (currently amended): The method of Claim 123 wherein R¹⁵ is selected from H. eptionally substituted pyrrolidinyl, optionally substituted piperazinyl, optionally substituted piperidinyl, merpholinyl, 1,2,3,6 tetrahydro pyridinyl, (optionally substituted pyrrelidinyl) C1-G2-alkyl. (optionally substituted piperidinyl) G1-G2-alkyl. (optionally substituted piperazinyl) G1-C2-alkyl, morpholinyl-C1-C2-alkyl, C1-C4-alkylamino-C1-C4-alkyl, C1-C4-hydroxyalkylamino, (optionally substituted pyrrolidinyl) C. C. alkylamine, (optionally substituted piperidinyl) C. G2-alkylamino, (optionally substituted piperazinyl) G1-G2-alkylamino, morpholinyl G1-G2alkylamine, optionally substituted pyrrolidinyl-C₁-C₄-alkoxy, optionally substituted azetidinyl-C₁-C₄-alkoxy, tetrahydrofuryl-C₁-C₄-alkoxy, optionally substituted piperidinyl-C₁-C₄-alkoxy, C₁-C₄-alkylamino-C₁-C₄-alkoxy, tetrahydrofuryloxy, optionally substituted piperidinyloxy, and optionally substituted phenoxy-C₁-C₂-alkylaminoearbonyl and C₁-C₄-alkylaminothiocarbonyl; wherein R¹⁶ is selected from H, 5-6-membered nitrogen containing heterocyclylearbonyl, C₁-C₄- alkylaminocarbonyl, and C₁-C₄-alkylaminomethyl, and 5-6-membered nitrogen containing heteroevelylmethyl; and wherein R¹⁷ is selected from halo, C₁-C₂-alkyl, eptionally substituted 5 6-membered hetereary sulfernyl-C1-C2-alkyl, optionally substituted phonoxy; and C3-C6cycloalkyl-C2-C4-alkynyl+ and pharmaceutically acceptable derivatives thereof.

Claim 125 (currently amended): The method of Claim 124 wherein R¹⁵ is selected from H₅ tetrahydro-furanyloxy, 1-methylpyrrolidin-2-ylmethoxy, 2-pyrrolidinylmethoxy, 3-

pyrrolidinylmethoxy, 1-Boc-pyrrolidin-2-ylmethoxy, 4-piperidinylmethoxy, 1-Boc-piperidin-4-ylmethoxy, 1-Boc-piperidin-4-ylethoxy, 1-methyl-piperidin-4-ylmethoxy, 1-Boc-azetidin-3-ylmethoxy, 1-methyl-azetidin-3-ylmethoxy, 3-azetidinylmethoxy, 1-methyl-piperidin-4-yloxy, phenyloxy, phenyloxy, 4-(pyrrolidin-1-ylmethyl)phenoxy, and dimethylaminoethoxy, 1-piperidinylmethyl, 1-(piperidin-1-yl)othyl, 3-methylpiperidin-1-ylmethyl, 1-pyrrolidinylmethyl, 2,2,6,6 tetramethylpiperidin-1-ylmethyl, 2,6 dimethylpiperidin-1-ylmethyl, dimethylaminomethyl, diethylaminomethyl, diethylaminomethyl, diethylaminomethyl, diethylaminomethyl, 2-thienylsulfonylmethyl, hydroxypropylamino, 4-ethyl-piperidin-1-yl, 4-(2-pyridyl)piperidin-1-yl, 1-methyl-1,2,3,6-tetrahydro-pyridin-4-yl, 1,2,3,6-tetrahydro-pyridin-4-yl, 4-(2-pyrazinyl)piperidin-1-yl, 1-methyl-1,2,3,6-tetrahydro-pyridin-4-yl, piperidinylmethyl, and 1-Boc-1,2,3,6-tetrahydro-pyridin-4-yl, wherein R¹⁶ is selected from H, 1-piperidinylearbonyl, diethylaminocarbonyl, and diethylaminomethyl, piperidinylmethyl, and pharmacoutically acceptable derivatives thereof.

Claim 126 (currently amended): The method of Claim 125, wherein R¹⁷ is chloro or bromo; and pharmaceutically acceptable derivatives thereof.

Claim 127 (currently amended): The method of Claim 123, wherein R¹⁵ is selected from H, optionally substituted piperazinyl, optionally substituted piperazinyl, optionally substituted piperazinyl, optionally substituted piperazinyl, (optionally substituted piperazinyl) C₄-C₂-alkyl, (optionally substituted piperazinyl) C₄-C₂-alkyl, morpholinyl C₄-C₂-alkyl, (optionally substituted piperazinyl) C₄-C₂-alkylamino, (optionally substituted piperazinyl) C₄-C₂-alkylamino, morpholinyl C₄-C₂-alkylamino, (optionally substituted piperazinyl) C₄-C₄-alkylamino, morpholinyl C₄-C₂-alkylamino, C₄-C₄-alkylamino C₄-C₄-alkyl₄-C₄-hydroxyalkylamino, optionally substituted pyrrolidinyl-C₁-C₄-alkoxy, optionally substituted azetidinyl-C₁-C₄-alkoxy, tetrahydrofuryl-C₁-C₄-alkoxy, optionally substituted piperidinyl-C₁-C₄-alkoxy, tetrahydrofuryloxy, optionally substituted piperidinyloxy, and optionally substituted phenoxy; C₄-C₄-alkylaminocarbonyl and C₄-C₄-alkylaminocarbonyl and C₄-C₄-alkylaminocarbonyl, C₁-C₄-alkylaminocarbonyl, and C₁-C₄-alkylaminomethyl; and 5-6-

membered-nitrogen containing heterocyolylmethyl; and wherein R¹⁷ is selected from C₃-C₆-cycloalkyl and phenyl-optionally substituted with one or two substituents selected from halo;

C₄-C₄-alkylamino, amino, nitro, C₄-C₄-alkoxy, C₄-C₂-halcalkyl, hydroxy, C₄-C₄-alkylthio, C₄-C₄-alkylcarbonylamino, (optionally substituted phonyl)sulfonylamino, eyano, C₄-C₂-halcalkoxy, 5- or 6-membered N-containing heterocyclyl, aminosulfonyl, (6-membered N-containing-heterocyclyl)sulfonyl, C₄-C₂-halcalkylcarbonylaminosulfonyl and (optionally substituted-phonyl)aminosulfonyl;

and pharmacoutically acceptable derivatives thereof.

Claim 128 (currently amended): The method of Claim 127, wherein R¹⁵ is selected from H₂ tetrahydro-furanyloxy, 1-methylpyrrolidin-2-ylmethoxy, 2-pyrrolidinylmethoxy, 3pyrrolidinylmethoxy, 1-Boc-pyrrolidin-2-ylmethoxy, 4-piperidinylmethoxy, 1-Boc-piperidin-4vlmethoxy, 1-Boc-piperidin-4-ylethoxy, piperidin-4-ylethoxy, 1-methyl-piperidin-4-ylmethoxy, 1-Boc-azetidin-3-ylmethoxy, 1-methyl-azetidin-3-ylmethoxy, 3-azetidinylmethoxy, 1-methylpiperidin-4-yloxy, phenyloxy; phenoxy 4-(pyrrolidin-1-ylmethyl)phenoxy, and dimethylaminoethoxy, 1-piperidinylmethyl, 1-(piperidin-1-yl)ethyl, 3-methylpiperidin-1ylmethyl, 1-pyrrolidinylmethyl, 2,2,6,6-tetramethylpiperidin-1-ylmethyl, 2,6-dimethylpiperidin-1-vlmethyl-dimethylaminomethyl, diethylaminomethyl, diethylaminothiocarbonyl, diethylaminocarbonyl, N. Boe-N. isopropylaminomethyl, isopropylaminomethyl, 2thionylsulfonylmethyl, hydroxypropylamino, 4-ethyl-piperidin-1-yl, 4-(2-pyridyl)piperidin-1-yl, 1 methylpiperidin 4 yl, 4 (2 pyrazinyl)piperidin 1 yl, 1 methyl 1,2,3,6 tetrahydro pyridin 4 yl, 1.2.3.6 tetrahydro pyridin 4 yl. and 1 Boc 1.2.3.6 tetrahydro pyridin 4 yl; wherein R¹⁶ is selected from H. 1-piperidinylearbonyl. diethylaminocarbonyl, and diethylaminomethyl. 1-piperidinylmethyl: and wherein R¹⁷ is selected from cyclopropyl and phenyl-eptionally substituted with aminosulfonyl; and pharmacoutically acceptable derivatives thereof.

Claim 129 (currently amended): The method of Claim 128, wherein R¹⁷ is unsubstituted phenyls and pharmaceutically acceptable derivatives thereof.

Claim 130 (currently amended): The method of Claim 123, wherein R¹⁵ is selected from H, optionally substituted piperazinyl, optionally substituted piperazinyl, optionally substituted piperazinyl, optionally substituted piperazinyl, optionally substituted pyrrolidinyl) G₁-

G₂-alkyl, (optionally substituted piperidinyl) G₄-C₂-alkyl, (optionally substituted piperazinyl) G₄-C₂-alkyl, morpholinyl G₄-C₂-alkyl, (optionally substituted pyrrolidinyl) G₄-C₂-alkylamino, (optionally substituted piperazinyl) G₄-C₂-alkylamino, (optionally substituted piperazinyl) G₄-C₂-alkylamino, morpholinyl G₄-C₂-alkylamino, G₄-G₄-alkylamino G₄-G₄-alkyl, G₄-G₄-alkylamino, optionally substituted pyrrolidinyl-C₁-C₄-alkoxy, optionally substituted azetidinyl-C₁-C₄-alkoxy, tetrahydrofuryl-C₁-C₄-alkoxy, optionally substituted piperidinyl-C₁-C₄-alkoxy, C₁-C₄-alkylamino-C₁-C₄-alkoxy, tetrahydrofuryloxy, optionally substituted piperidinyloxy, and optionally substituted phenoxy, G₄-G₄-alkylaminocarbonyl and G₄-G₄-alkylaminothiocarbonyl, wherein R¹⁶ is selected from H, 5-6-membered nitrogen-containing heterocyclylaminopholyl, and C₁-C₄-alkylaminomethyl, and 5-6-membered nitrogen-containing heterocyclylmethyl; and wherein R¹⁷ is selected from optionally substituted indazelyl, optionally substituted indelyl, unsubstituted 5-membered exygen or sulfur containing heterocyclyl, and 6-membered nitrogen-containing heterocyclyl substituted with one or more substituents independently colected from pyridyl, phenyl,

C₄-C₄-alkyl, C₄-C₂-haloalkyl, C₁-C₂ alkoxy, amino, halo, piperidinyl, morpholinyl, C₄-C₂ alkylamino-C₄-C₂ alkylamino-C₄-C₄-alkylamino-C₄-C₄-alkylamino-C₄-C₄-alkylamino-C₄-C₄-alkylamino-C₄-C₄-alkylamino-C₄-C₄-alkylamino-C₄-C₄-haloalkylearbonylamino, morpholinyl-C₄-C₄-alkylamino, N,N-di-C₄-C₂-alkylamino-C₄-C₄-Alkylamino-C₄-C₄-Alkylamino-C₄-C₄-Alkylamino-C₄-C₄-Alkylamino-C₄-C₄-Alkylamino-C₄-C₄-Alkylamino-C₄-C₄-Alkylamino-C₄-C₄-Alkylamino-C₄-C₄-Alkylamino-C₄-C

and pharmaceutically acceptable derivatives thereof.

Claim 131 (currently amended): The method of Claim 130, wherein R¹⁵ is selected from H, tetrahydro-furanyloxy, 1-methylpyrrolidin-2-ylmethoxy, 2-pyrrolidinylmethoxy, 3-pyrrolidinylmethoxy, 1-Boc-pyrrolidin-2-ylmethoxy, 4-piperidinylmethoxy, 1-Boc-piperidin-4-ylethoxy, piperidin-4-ylethoxy, 1-methyl-piperidin-4-ylmethoxy, 1-Boc-azetidin-3-ylmethoxy, 1-methyl-azetidin-3-ylmethoxy, 3-azetidinylmethoxy, 1-methyl-piperidin-4-yloxy, phenyloxy, 4-(pyrrolidin-1-ylmethyl)phenoxy, and dimethylaminoethoxy; 1-piperidinylmethyl, 1-(piperidin-1-yl)ethyl, 3-methylpiperidin-1-ylmethyl, 1-pyrrolidinylmethyl, 2,2,6,6-totramethylpiperidin-1-ylmethyl, 2,6-dimethylpiperidin-1-ylmethyl, diethylaminoethyl, diethylaminoethylaminoethyl, diethylaminoethylaminoethylaminoethylaminoethylaminoethylaminoethyla

Boe N isopropylaminomethyl, isopropylaminomethyl, 2 thionylsulfonylmethyl, hydroxypropylamino, 4 ethyl piperidin 1 yl, 4 (2 pyridyl)piperidin 1 yl, 1 methyl piperidin 1 yl, 4 (2 pyridyl)piperidin 1 yl, 1 methyl 1,2,3,6 tetrahydro pyridin 4 yl, 1,2,3,6 tetrahydro pyridin 4 yl, 1,2,3,6 tetrahydro pyridin 4 yl; wherein R¹⁶ is selected from H, 1 piperidinylearbonyl, diethylaminocarbonyl, and diethylaminomethyl, 1 piperidinylmethyl; and wherein R¹⁷ is selected from 5 indazolyl, 1 Boe indol 5 yl, unsubstituted thienyl, 5 tert butylexazol 2 yl and 4 pyridyl substituted with one or more substituents independently selected from methoxy and chloro; and pharmaccutically acceptable derivatives thereof.

Claim 132 (currently amended): The method of Claim 130 123, wherein R¹⁷ is 4-pyridyl; and pharmaceutically acceptable derivatives thereof.

Claim 133 (currently amended): The method of Claim 123, wherein the compound is and pharmaceutically acceptable derivatives thereof selected from:

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1-[6-(3-Methyl-piperidin-1-ylmethyl)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;
1-[4 (Piperidine-1 earbonyl) pyridin-2-yl] 3 (2-pyridin-4-yl-thiazel-4-yl) urea;
1-(2-Chloro thiazol 4-yl) 3-[4 (piperidine-1-earbonyl) pyridin-2-yl] urea;
N.N. Diethyl 2 [3 (2 pyridin 4 yl-thiazol 4-yl) urcido] isonicotinamide;
N.N. Diethyl 2-[3-(2-phonyl thiazel-4-yl) ureido] isonicotinamide;
2-[3-(2-Brome-thiazel-4-yl)-ureide]-N;N-diethyl-isenicetinamide;
1-(4-Diethylaminomethyl-pyridin-2-yl)-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
1-16 (2.6 Dimethyl piperidin 1-ylmethyl) pyridin 2-yl 3 (2-pyridin 4-yl thiazol 4-yl) urea;
4-[6 (1 Piperidin 1 yl ethyl) pyridin 2 yl] 3 (2 pyridin 4 yl thiazol 4 yl) urea;
2 ({6 [3 (2 Pyridin 4 yl thiazol 4 yl) urcido} pyridin 2 ylamino}-methyl) piperidine 1-
     earboxylic acid tert butyl ester:
1 (6 [(Piperidin 2 ylmothyl) amino] pyridin 2 yl}-3 (2 pyridin 4 yl thiazol 4 yl) urea;
(S) 1 [6 (3 Methyl piperidin 1 ylmethyl) pyridin 2 yl] 3 (2 pyridin 4 yl thiazol 4 yl) urea;
(R) 1 [6 (3 Mothyl piperidin 1 ylmothyl) pyridin 2 yl] 3 (2 pyridin 4 yl thiazel 4 yl) urea;
1-(2 Chloro thiazol 4 vl) 3-(6 piperidin 1-ylmethyl-pyridin 2 yl) urea;
1-(2-Bromo-thiazol-4-yl)-3-[6-(2-piperidin-4-yl-ethoxy)-pyridin-2-yl]-urea;
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1-(2-Chloro-thiazol-4-yl)-3-[6-(2-piperidin-4-yl-ethoxy)-pyridin-2-yl]-urea;
1-[6-(Azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-bromo-thiazol-4-yl)-urea;
1-[6-(Azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-chloro-thiazol-4-yl)-urea;
1-(2-Bromo-thiazol-4-yl)-3-[6-(piperidin-4-ylmethoxy)-pyridin-2-yl]-urea;
1-(2-Chloro-thiazol-4-yl)-3-[6-(piperidin-4-ylmethoxy)-pyridin-2-yl]-urea;
tert-Butyl 3-{6-[3-(2-pyridin-4-yl-thiazol-4-yl)-ureido]-pyridin-2-yloxymethyl}-pyrrolidine-1-
     carboxylate;
1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(pyrrolidin-3-ylmethoxy)-pyridin-2-yl]-urea;
1-(2-Cyclopropyl-thiazol-4-yl)-3-[6-(2-piperidin-4-yl-ethoxy)-pyridin-2-yl]-urea;
1-[6 (Isopropylamino-methyl) pyridin-2-yl]-3 (2-pyridin-4-yl-thiazol-4-yl) urea;
1-[6 (Isopropylamino-methyl) pyridin-2-yl]-3-(2-phonyl-thiazol-4-yl) urea;
1-(2-Bromo-thiazol-4-yl)-3-[6-(isopropylamino-methyl)-pyridin-2-yl]-urea;
1-(2-Bromo-thiazol-4-yl)-3-[6-(1-methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-urea;
1-(2-Chloro-thiazol-4-yl)-3-[6-(1-methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-urea;
1-(2-phenylthiazol-4-yl)-3-(6-p-pyrrolidin-1-ylmethylphenoxypyridin-2-yl)urea;
1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(tetrahydro-furan-3-yloxy)-pyridin-2-yl]-urea;
1-12-(1H-Indezel-5-vl) thiezel 4-vl] 3-(6-piperidin 1-vlmethyl pyridin 2-yl) urea;
1-(1' Methyl-1',2',3',6' tetrahydro-[2,4']bipyridinyl-6-yl) 3-(2-pyridin-4-yl-thiazol-4-yl) uroa;
1 (2 Bromo thizol 4 yl) 3 (1' methyl-1':2':3':6' tetrahydro [2:4']bipyridinyl 6 yl) urca;
1-(1'-Methyl-1',2',3',6'-tetrahydro-2[2,4]bipyridinyl-6-yl)-3-(2-phenyl-thiazel-4-yl)-urea;
1-[6 (3 Hydroxy propylamino) pyridin 2 yl] 3 (2 pyridin 4 yl-thizol 4 yl) urea;
1-(2-Bromo thiazol-4-yl) 3-[6(3-hydroxy-propylamino) pyridin-2-yl]-uroa;
1-(1' Methyl-1',2',3',4',5',6' hexahydro-[2,4']bipydrinyl-6-yl) 3-(2-pyridin-4-yl-thiazol-4-yl)
     urca:
1 (1' Methyl 1',2',3',4',5',6' hexahydro [2,4']bipyridinyl 6 yl) 3 (2 phonyl thiazol 4 yl) uroa;
6 [3 (2 Pyridin 4 yl thizol 4 yl) ureide] 3',6' dihydro 2'H [2,4]bipyridinyl 1' carboxylio acid
     tert-butylester;
1-(2-Pyridin 4 yl thiazol 4 yl) 3-(1',2',3',6'-tetrahydro-[2,4']bipyridinyl-6-yl) urea;
1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(tetrahydro-furan-3-ylmethoxy)-pyridin-2-yl]-urea;
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- 2-[6-[3-(2-Pyridin-4-yl-thiazol-4-yl)-ureido]-pyridin-2-yloxymethyl]-pyrrolidine-1-carboxylic acid tert-butyl ester;
- 1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-urea;
- 6 [3 (2 Pyridin 4 yl thiazol 4 yl) urcido} pyridine 2 carbothioic acid diethylamide;
- 1-(2-Bromo-thiazol-4-yl)-3-[6-(3-methyl-piperidin-1-ylmethyl)-pyridin-2-yl]-urea;
- 1-(2-Chloro-thiazol-4-yl)-3-[6-(3-methyl-piperidin-1-ylmethyl)-pyridin-2-yl]-urea;
- 1-(2 Phenyl thiazol 4-yl)-3-[4 (piperidino-1-carbonyl)-pyridin-2-yl]-urea;
- 1 (2 Brome thiazel 4 yl) 3 [4 (piperidine 1 earbonyl) pyridin 2 yl] urea;
- 1-[2-(2-Methoxy-pyridin-4-yl)-thiazol-4-yl]-3-(6-phenoxy-pyridin-2-yl)-urea;
- 1-[2-(2-Methoxy-pyridin-4-yl)-thiazol-4-yl]-3-[6-(1-methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-urea;
- 1-[6-(2-Dimethylamino-ethoxy)-pyridin-2-yl]-3-[2-(2-methoxy-pyridin-4-yl)-thiazol-4-yl]-urea;
- 1-[6-(1-Methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-(2-phenylthiazol 4-yl)-3-(6-pyrrolidin-1-ylmethyl-pyridin-2-yl)urca;
- 1-(6-Diethylaminomethylpyridin 2 yl)-3-(2-phenylthiazol-4-yl)urea;
- (S)-1-[6-(1-Methylpyrrolidin-2-ylmethoxy)pyridin-2-yl]-3-(2-phenylthiazol-4-yl)urea;
- 1-[6-(2-Piperidin-4-yl-ethoxy)pyridin-2-yl]-3-[2-phenylthiazol-4-yl]urea;
- 1-[6 (4-Ethylpiporazin 1-yl) pyridin 2-yl] 3 (2-phenylthiazol 4-yl)urca;

Diethyl 6-[3-(2-phenylthiazol 4-yl)uroido]-pyridino-2-earboxamide;

- 1-(2-Pyridin-4-yl-thiazol-4-yl)-3-(6-p-pyrrolidin-1-ylmethylphenoxypyridin-2-yl)urea;
- 1-(2-Bromothiazol-4-yl)-3-(6-p-pyrrolidin-1-ylmethylphenoxypyridin-2-yl)urea;
- 1-[6-(Piperidin-4-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-piperidin-4-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-piperidin-4-yloxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-[6-(Azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;
- 1-(2-Phenyl-thiazol-4-yl)-3-[6-(piperidin-4-ylmethoxy)-pyridin-2-yl]-urea;
- 1-[6-(1-Methyl-piperidin-4-ylmethoxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-piperidin-4-yloxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;
- 1-[6-(2-Piperidin-4-yl-ethoxy)-pyridin-2-yl]-3-(2-thiophen-2-yl-thiazol-4-yl)-urea; and

- 1-[6-(1-Methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-3-[2-(thiophene-2-sulfonylmethyl)-thiazol-4-yl]-urea;
- 1-[2 (2 Methoxy pyridin 4 yl) thiazel 4 yl] 3 (6 piperdin-1-ylmethyl-pyridin 2 yl) urea; and [2-(2 Chloro-pyridin 4 yl) thiazel 4 yl] 3 (6 piperdin-1-ylmethyl-pyridin-2 yl) urea; and pharmaceutically acceptable salts thereof.

Claim 134 (currently amended): The method of Claim 123, wherein the compound is and pharmaceutically acceptable derivatives thereof selected from:

- 1-[6-(3-Methyl-piperidin-1-ylmothyl)-pyridin-2-yl]-3-(2-phonyl-thiazol-4-yl)-urea;
- 1-[4-(Piperidino-1-earbonyl) pyridin 2-yl]-3-(2-pyridin 4-yl-thiazol-4-yl) urca;
- N.N-Diethyl-2-[3-(2-pyridin-4-yl-thiazol-4-yl)-uroido]-isonicotinamide;
- 1-(4-Diethylaminomethyl-pyridin-2-yl) 3-(2-pyridin-4-yl-thiazol-4-yl) urea;
- 1-[6-(2,6-Dimothyl-piperidin-1-ylmothyl)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-[6 (1 Piperidin 1 yl ethyl) pyridin 2 yl] 3 (2 pyridin 4 yl thiazel 4 yl) urea;
- 2 ({6 [3 (2 Pyridin 4 yl thiazel 4 yl) ureide] pyridin-2 ylamine} methyl) piperidine 1 carboxylic acid tert-butyl ester:
- 1 (6 (Piperidin 2 ylmethyl)-amino pyridin 2 yl) 3 (2 pyridin 4 yl thiazol 4 yl) urea;
- (S) 1-[6 (3 Methyl-piperidin-1-ylmethyl) pyridin-2-yl] 3 (2 pyridin-4-yl-thiazol-4-yl) urea;
- (R) 1-[6 (3 Methyl-piperidin-1-ylmethyl) pyridin-2-yl] 3 (2 pyridin-4-yl-thiazol-4-yl) urea;
- 1-(2-Chloro-thiazol-4-yl)-3-(6-piperidin-1-ylmethyl-pyridin-2-yl)-urea;
- 1-(2-Bromo-thiazol-4-yl)-3-[6-(2-piperidin-4-yl-ethoxy)-pyridin-2-yl]-urea;
- 1-(2-Chloro-thiazol-4-yl)-3-[6-(2-piperidin-4-yl-ethoxy)-pyridin-2-yl]-urea;
- 1-[6-(Azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-bromo-thiazol-4-yl)-urea;
- 1-[6-(Azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-chloro-thiazol-4-yl)-urea;
- 1-(2-Bromo-thiazol-4-yl)-3-[6-(piperidin-4-ylmethoxy)-pyridin-2-yl]-urea;
- 1-(2-Chloro-thiazol-4-yl)-3-[6-(piperidin-4-ylmethoxy)-pyridin-2-yl]-urea;
- 3-(4-{3-[6-(1-Methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-ureido}-thiazol-2-yl)-benzenesulfonamide;
- *tert*-Butyl 3-{6-[3-(2-pyridin-4-yl-thiazol-4-yl)-ureido]-pyridin-2-yloxymethyl}-pyrrolidine-1-carboxylate;
- 1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(pyrrolidin-3-ylmethoxy)-pyridin-2-yl]-urea;

- 1-(2-Cyclopropyl-thiazol-4-yl)-3-[6-(2-piperidin-4-yl-ethoxy)-pyridin-2-yl]-urea;
- Isopropyl (6 [3 (2 pyridin 4 yl-thiazol-4 yl) urcido] pyridin-2-ylmethyl) carbamic acid tertbutyl ester;
- 1-[6 (Isopropylamino-methyl) pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl) urea;
- Isopropyl (6 [3 (2 phenyl-thiazol 4 yl) urcido]-pyridin-2-ylmethyl)-earbamic acid-tert-butyl
- 1 [6 (Isopropylamino methyl) pyridin 2 yl] 3 (2 phonyl thiazol 4 yl) urea;
- 1-(2-Bromo-thiazol-4-yl)-3-[6-(1-methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-urea;
- 1-(2-Chloro-thiazol-4-yl)-3-[6-(1-methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-urea;
- 1-(2-phenylthiazol-4-yl)-3-(6-p-pyrrolidin-1-ylmethylphenoxypyridin-2-yl)urea;
- 1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(tetrahydro-furan-3-yloxy)-pyridin-2-yl]-urea;
- 1 [2 (1H Indazol 5 yl) thiazol 4 yl] 3 (6 piperidin 1 ylmethyl pyridin 2 yl) urea;
- 1 (1' Methyl 1',2',3',6' tetrahydro [2,4']bipyridinyl 6 yl) 3 (2 pyridin 4 yl thiazol 4 yl) urea;
- 1-(2-Bromo-thizel-4-yl)-3-(12-methyl-12,22,32,62-tetrahydro-[2,42]bipyridinyl-6-yl)-urca;
- 1-(1'-Mothyl-1',2',3',6'-tetrahydro-2[2,4]bipyridinyl-6-yl)-3-(2-phonyl-thiazol-4-yl)-urea;
- 1-[6 (3 Hydroxy propylamino) pyridin-2-yl]-3 (2 pyridin-4-yl-thizol-4-yl) urea;
- 1 (2 Bromo thiazol 4 yl) 3 [6(3 hydroxy propylamino) pyridin-2-yl]-urea;
- 1-(1' Mothyl 1',2',3',4',5',6'-hexahydro [2,4']bipydrinyl 6-yl) 3-(2-pyridin-4-yl thiazol 4-yl) uroa;
- 1 (1' Methyl 1',2',3',4',5',6' hoxahydro [2,4']bipyridinyl 6 yl) 3 (2 phonyl thiazol 4 yl) urea;
- 6 [3 (2-Pyridin 4-yl-thizel 4-yl) urcide] 3',6'-dihydre 2'H-[2,4]bipyridinyl-1'-carboxylic-acid tert-butylester;
- 1-(2-Pyridin 4 yl-thiazol 4 yl) 3-(1',2',3',6'-tetrahydro-[2,4']bipyridinyl-6-yl)-uroa;
- 1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(tetrahydro-furan-3-ylmethoxy)-pyridin-2-yl]-urea;
- 2-[6-[3-(2-Pyridin-4-yl-thiazol-4-yl)-ureido]-pyridin-2-yloxymethyl]-pyrrolidine-1-carboxylic acid tert-butyl ester;
- 1-(2-Pyridin-4-yl-thiazol-4-yl)-3-[6-(pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-urea;
- 6-[3-(2-Pyridin 4-yl-thiazol 4-yl)-urcido]-pyridino-2-carbothioio acid-diethylamido;
- 1-(2-Bromo thiazol 4-yl)-3-[6-(3-methyl-piperidin-1-ylmethyl)-pyridin-2-yl]-urea;
- 1-(2-Chloro-thiazol-4-yl)-3-[6-(3-methyl-piperidin-1-ylmethyl)-pyridin-2-yl]-urea;
- 1-[2-(2-Methoxy-pyridin-4-yl)-thiazol-4-yl]-3-[6-(1-methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-urea;

- 1-[6-(2-Dimethylamino-ethoxy)-pyridin-2-yl]-3-[2-(2-methoxy-pyridin-4-yl)-thiazol-4-yl]-urea;
- 1-[6-(1-Methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-(2-phonylthiazol 4-yl)-3-(6-pyrrolidin-1-ylmethyl-pyridin-2-yl)urea;
- 1-(6-Diethylaminomethylpyridin 2 yl) 3-(2 phenylthiazol 4 yl)urea;
- (S)-1-[6-(1-Methylpyrrolidin-2-ylmethoxy)pyridin-2-yl]-3-(2-phenylthiazol-4-yl)urea;
- 1-[6-(2-Piperidin-4-yl-ethoxy)pyridin-2-yl]-3-[2-phenylthiazol-4-yl]urea;
- 1- [6 (4 Ethylpiporazin 1-yl) pyridin 2-yl] 3 (2 phonylthiazol 4-yl)urca;
- 1-(2-phenylthiazel-4-yl)-3-[6-(4-pyrimidin-2-yl-piperazin-1-yl)pyridin-2-yl]urea;

Diethyl 6-[3-(2-phonylthiazol-4-yl)urcido]-pyridine-2-carboxamide;

- 1-(2-Pyridin-4-yl-thiazol-4-yl)-3-(6-p-pyrrolidin-1-ylmethylphenoxypyridin-2-yl)urea;
- 1-(2-Bromothiazol-4-yl)-3-(6-p-pyrrolidin-1-ylmethylphenoxypyridin-2-yl)urea;
- 1-[6-(Piperidin-4-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-piperidin-4-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-piperidin-4-yloxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-pyridin-4-yl-thiazol-4-yl)-urea;
- 1-[6-(Azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-azetidin-3-ylmethoxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;
- 1-(2-Phenyl-thiazol-4-yl)-3-[6-(piperidin-4-ylmethoxy)-pyridin-2-yl]-urea;
- 1-[6-(1-Methyl-piperidin-4-ylmethoxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;
- 1-[6-(1-Methyl-piperidin-4-yloxy)-pyridin-2-yl]-3-(2-phenyl-thiazol-4-yl)-urea;
- 1-[6-(2-Piperidin-4-yl-ethoxy)-pyridin-2-yl]-3-(2-thiophen-2-yl-thiazol-4-yl)-urea; and
- 1-[6-(1-Methyl-pyrrolidin-2-ylmethoxy)-pyridin-2-yl]-3-[2-(thiophene-2-sulfonylmethyl)-thiazol-4-yl]-urea;
- 1-[2 (2 Mothoxy pyridin 4-yl) thiazol 4-yl] 3 (6 piperdin 1-ylmethyl pyridin 2-yl) urca; and
- [2 (2 Chloro pyridin 4 yl) thiazol-4 yl]-3 (6 piperdin-1-ylmethyl pyridin-2 yl) urca;

and pharmaceutically acceptable salts thereof.